

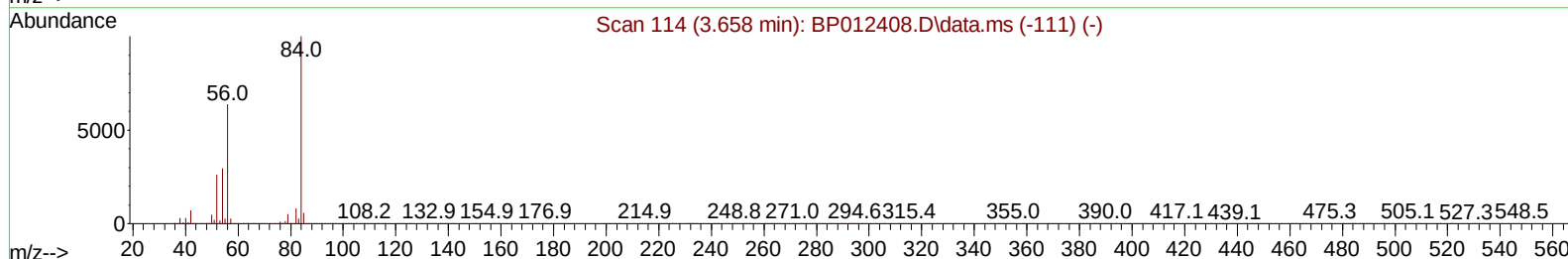
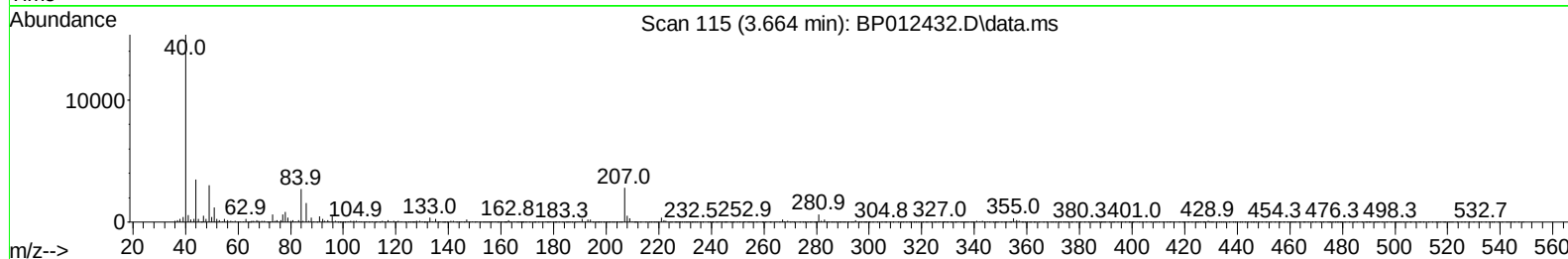
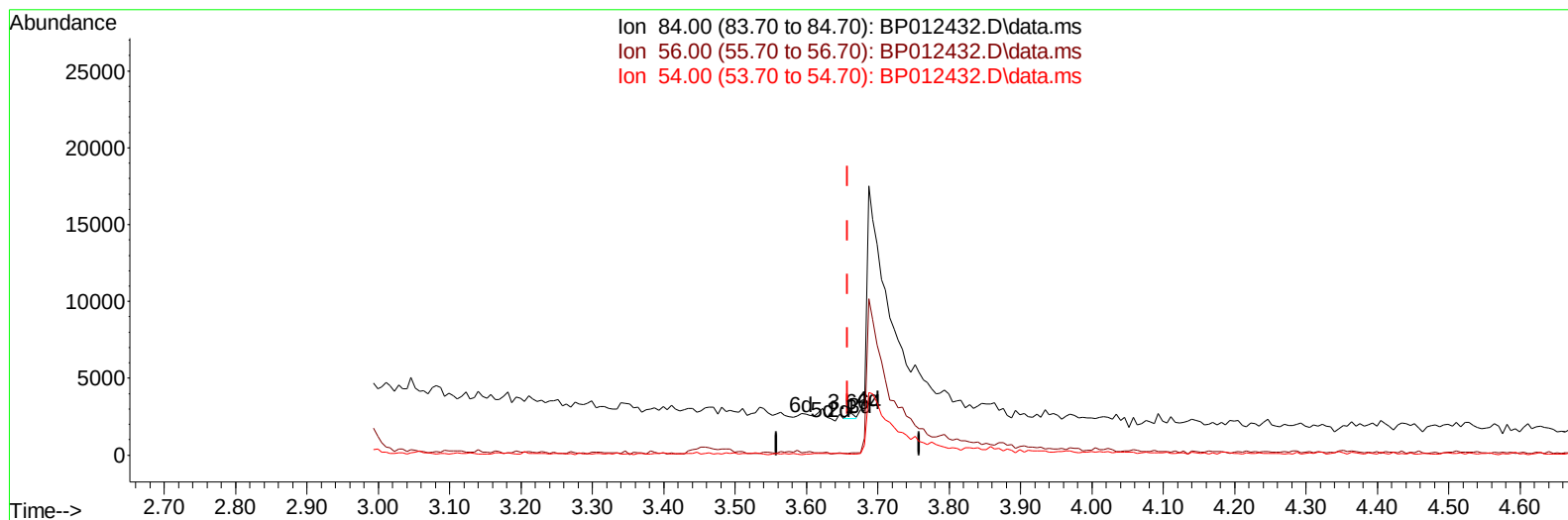
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
 Data File : BP012432.D
 Acq On : 01 Nov 2022 17:52
 Operator : CG/JU
 Sample : N5180-05DL 5X
 Misc :
 ALS Vial : 53 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0KX4DL

Manual IntegrationsAPPROVED

Quant Time: Nov 01 23:00:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 01 01:03:58 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/02/2022
 Supervised By :mohammad ahmed 11/02/2022



TIC: BP012432.D\data.ms

(4) Pyridine-d5 (S)

3.664min (+ 0.006) 0.01 ng/ul

response 223

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	5.13#
54.00	29.80	2.34#
0.00	0.00	0.00

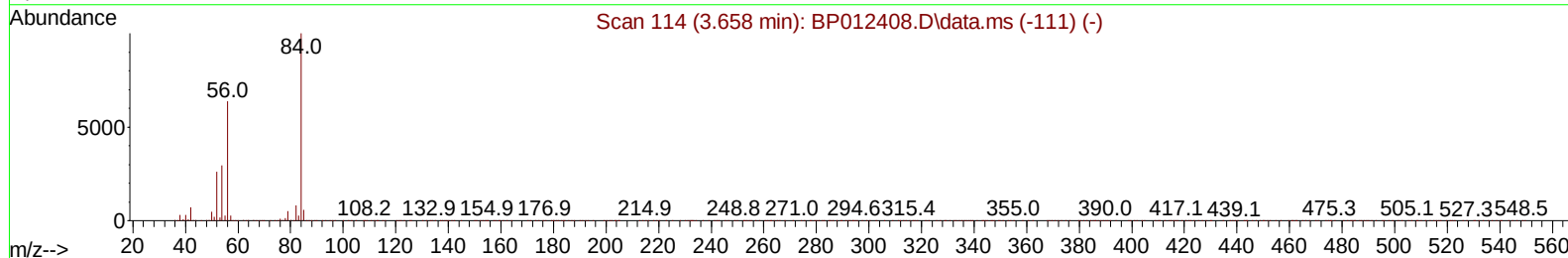
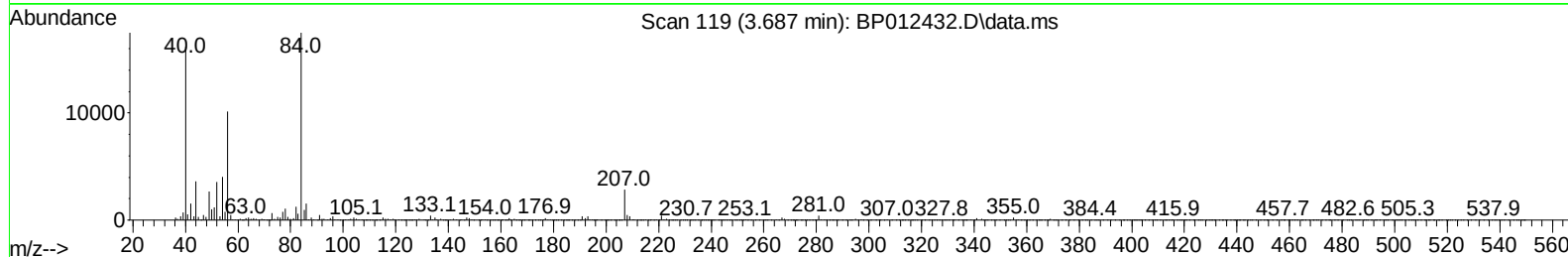
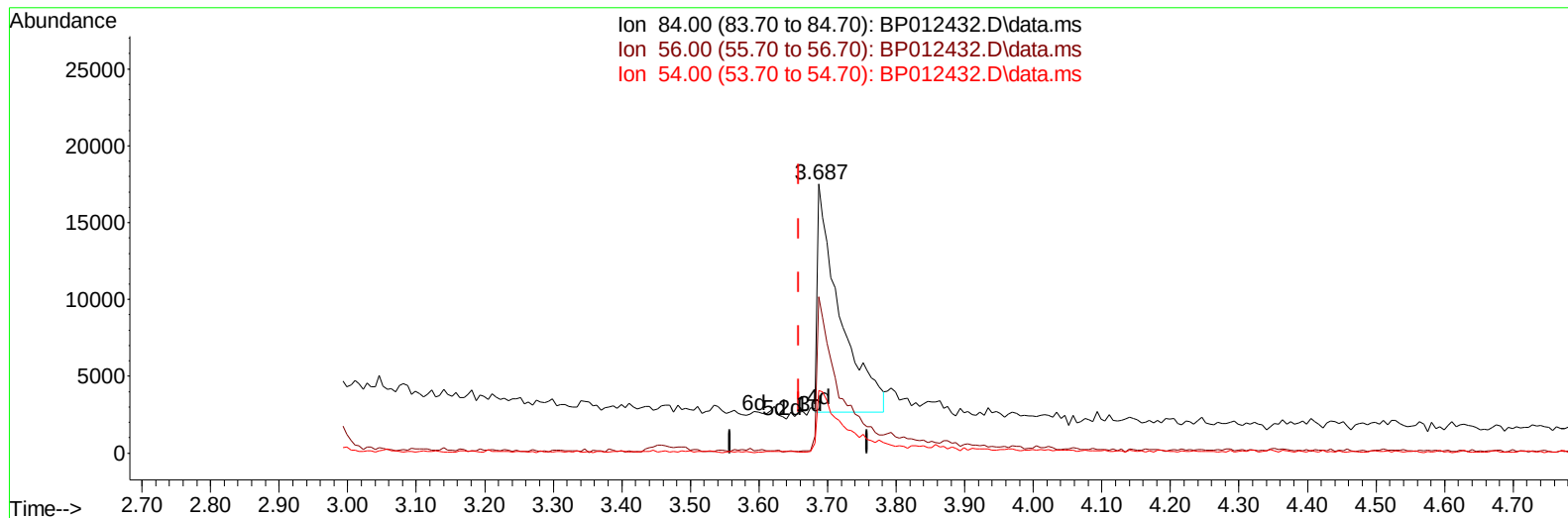
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(4) Pyridine-d5 (S)

3.687min (+ 0.029) 1.61 ng/ul m

response 33727

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	58.11
54.00	29.80	23.17#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	293615	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	1329697	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	878948	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1944044	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	2006295	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1786576	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	3.687	84	33727m	1.610	ng/ul	0.03
7) Phenol-d5	6.969	99	34358	1.305	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.134	67	119846	7.623	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	102835	5.347	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	66385	3.187	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	72473	8.405	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	64384	7.194	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	112528	5.802	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	139523	4.831	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	482243	8.015	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	517086	7.520	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.439	176	443634	8.612	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.304	188	706033	8.429	ng/ul	0.00
81) Pyrene-d10	19.545	212	862919	8.567	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.527	264	705827	8.086	ng/ul	-0.02
Target Compounds						
					Qvalue	
22) Nitrobenzene	9.005	77	433965	19.289	ng/ul	99
39) 1,2,4,5-Tetrachloroben...	12.622	216	284335	11.013	ng/ul	99
75) 1,2,3,4-Tetrachloroben...	13.234	216	701382	25.221	ng/uL	99

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

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