

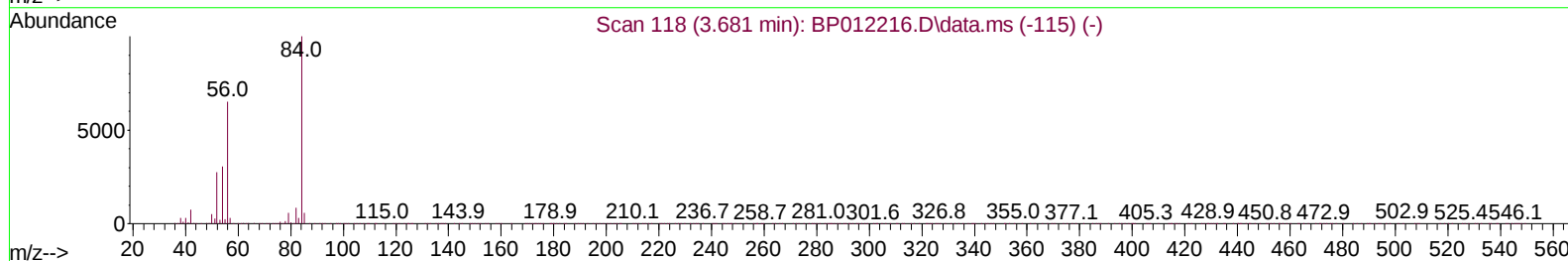
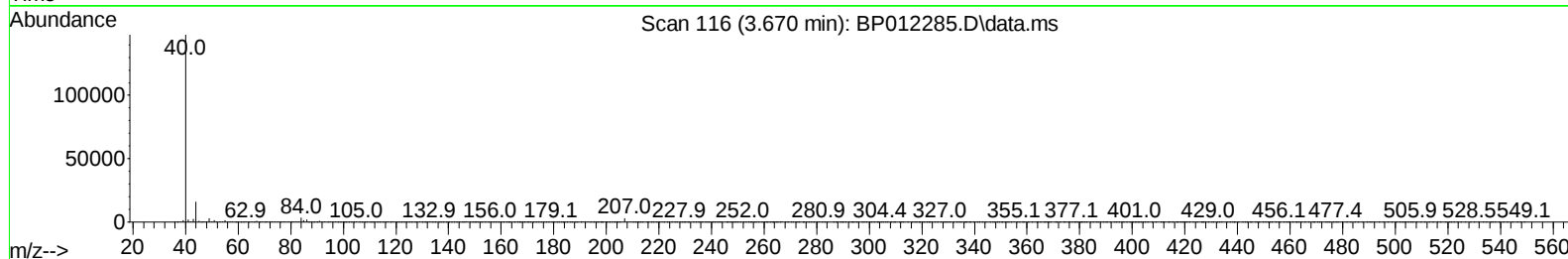
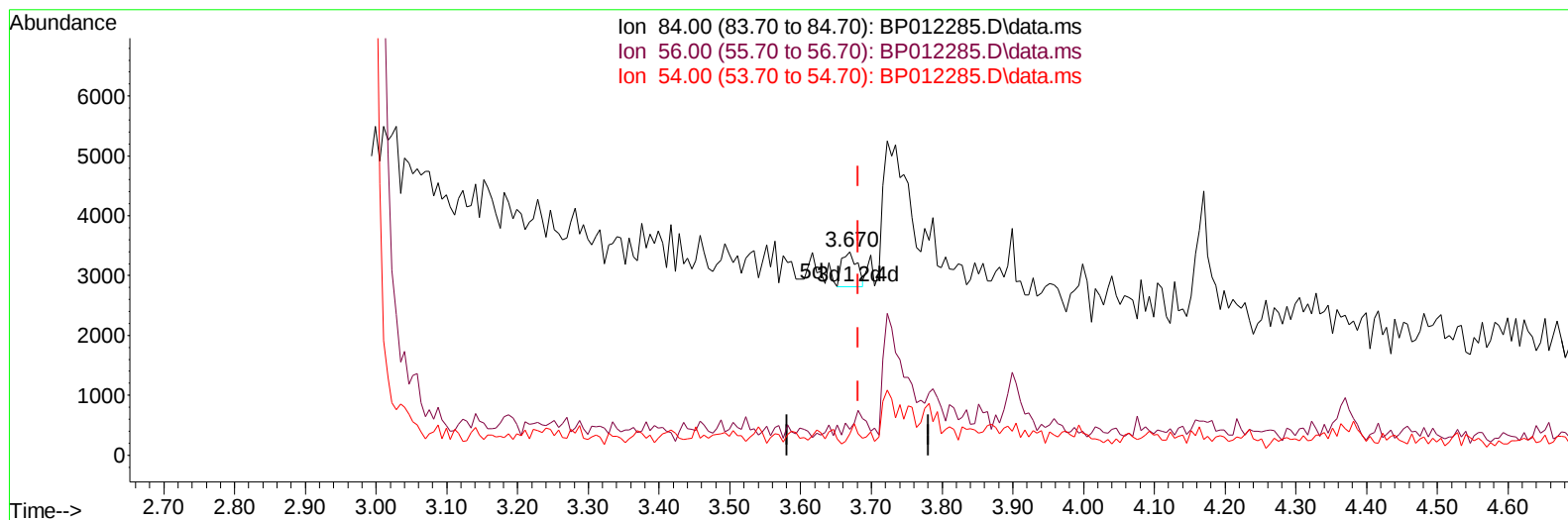
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012285.D
 Acq On : 24 Oct 2022 15:04
 Operator : CG/JU
 Sample : N5070-12
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BS2

Manual IntegrationsAPPROVED

Quant Time: Oct 25 05:53:33 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Oct 22 01:24:49 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/25/2022
 Supervised By :mohammad ahmed 10/27/2022



TIC: BP012285.D\data.ms

(4) Pyridine-d5 (S)

3.670min (-0.012) 0.03 ng/u1

response 860

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	12.71#
54.00	29.70	10.56#
0.00	0.00	0.00

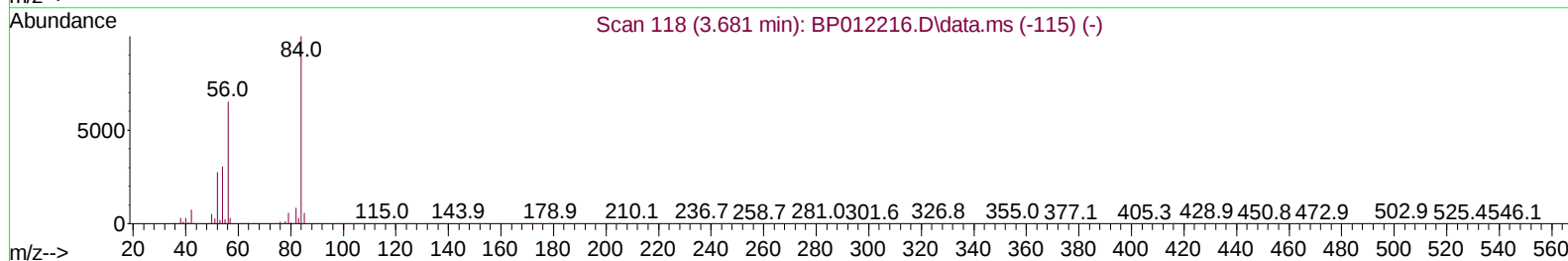
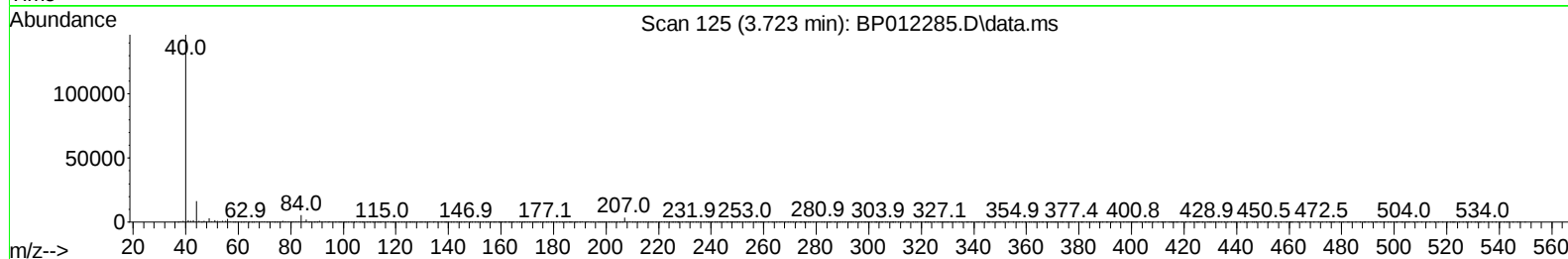
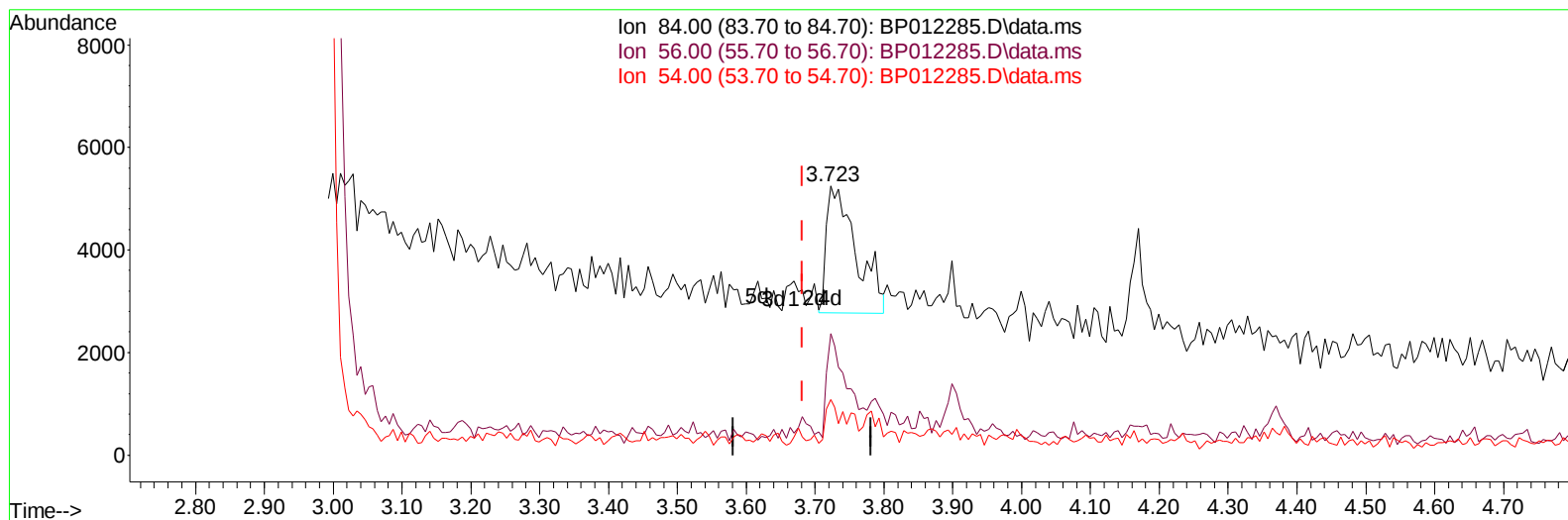
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TIC: BP012285.D\data.ms

(4) Pyridine-d5 (S)

3.723min (+ 0.041) 0.22 ng/u1 m

response 7390

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	45.19#
54.00	29.70	20.72#
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	7.816	152	465843	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	2017118	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	1200831	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	2157488	20.000	ng/ul	0.00
79) Chrysene-d12	21.316	240	1673934	20.000	ng/ul	0.00
88) Perylene-d12	23.716	264	1790562	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	39493	3.431	ng/uL	0.00
4) Pyridine-d5	3.723	84	7390m	0.223	ng/ul	0.04
7) Phenol-d5	6.987	99	947273	23.554	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.152	67	614003	25.174	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	769670	25.143	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	455316	14.567	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	425169	28.476	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	451642	28.036	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	732539	24.243	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	638490	14.872	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	2282861	27.329	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	2645956	27.751	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	261591	16.716	ng/ul	0.00
60) Fluorene-d10	15.463	176	1894522	26.363	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	161879	13.189	ng/ul	0.00
73) Anthracene-d10	17.328	188	2452711	25.936	ng/ul	0.00
81) Pyrene-d10	19.569	212	2626477	27.558	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.563	264	2406608	26.801	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	63337	1.505	ng/ul	95

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

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