

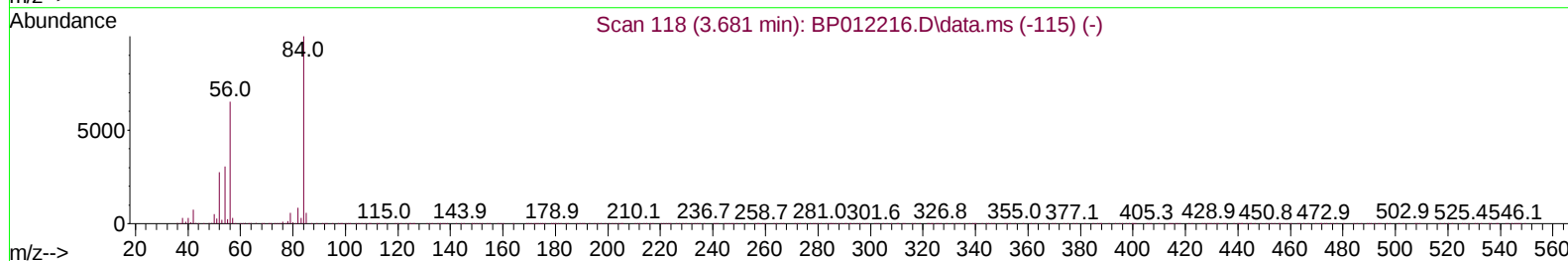
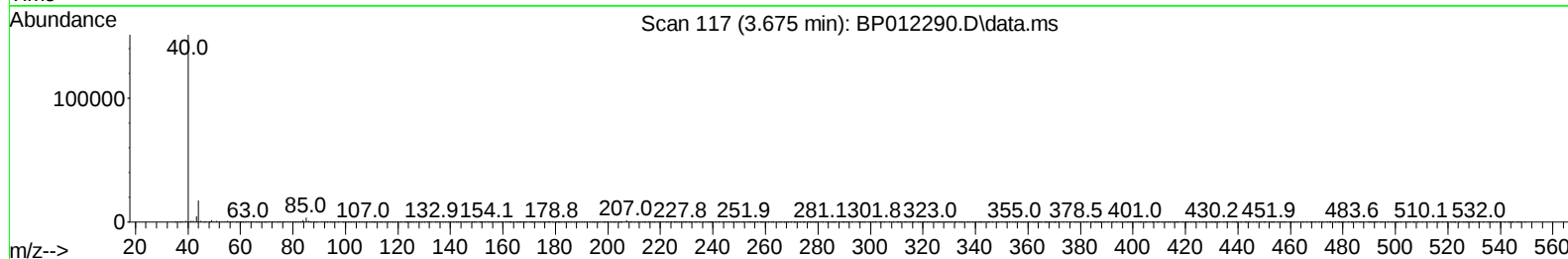
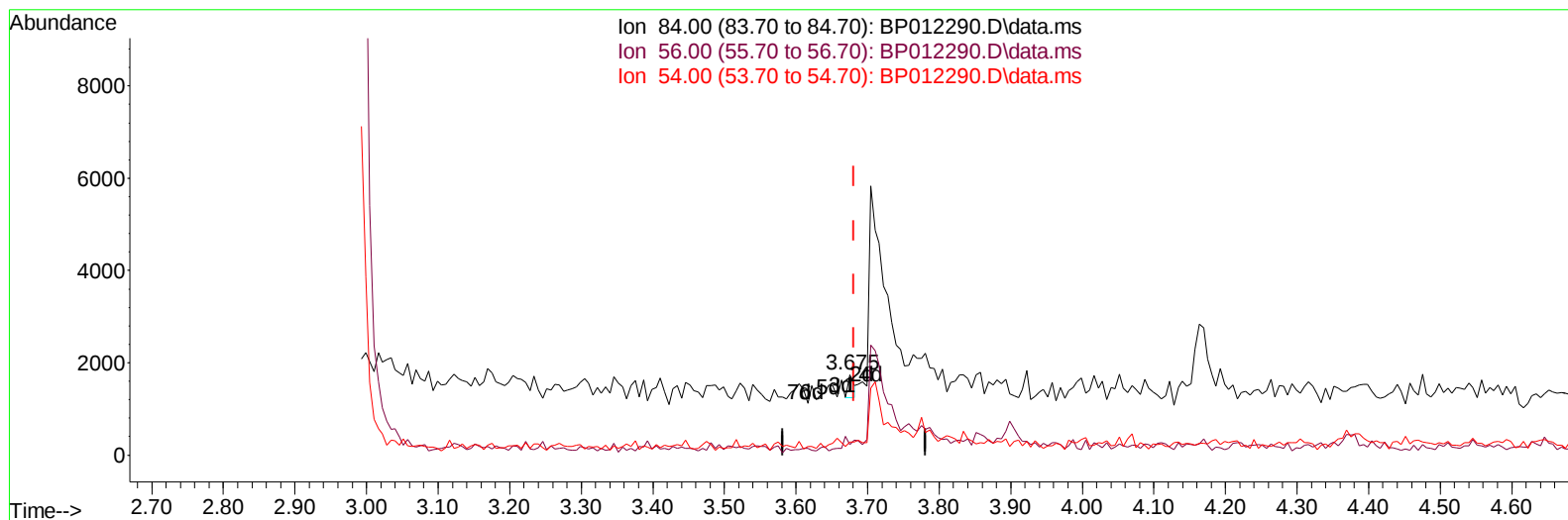
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102422\
 Data File : BP012290.D
 Acq On : 24 Oct 2022 17:56
 Operator : CG/JU
 Sample : N5070-11
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BR4

Manual IntegrationsAPPROVED

Quant Time: Oct 25 06:28:24 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: BP012290.D\data.ms

(4) Pyridine-d5 (S)

3.675min (-0.006) 0.01 ng/ul

response 264

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	12.86#
54.00	29.70	16.07#
0.00	0.00	0.00

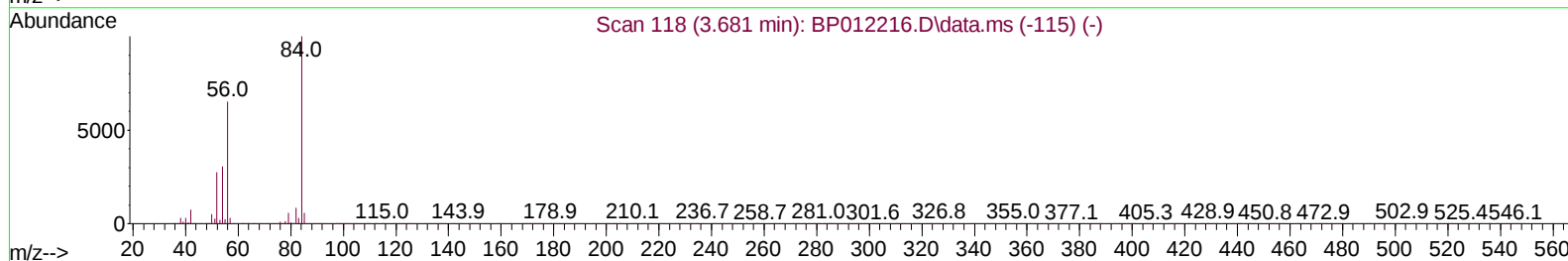
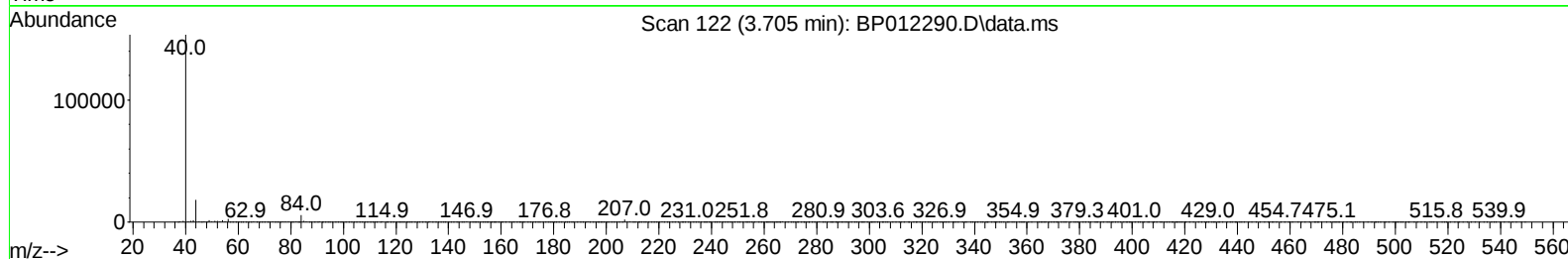
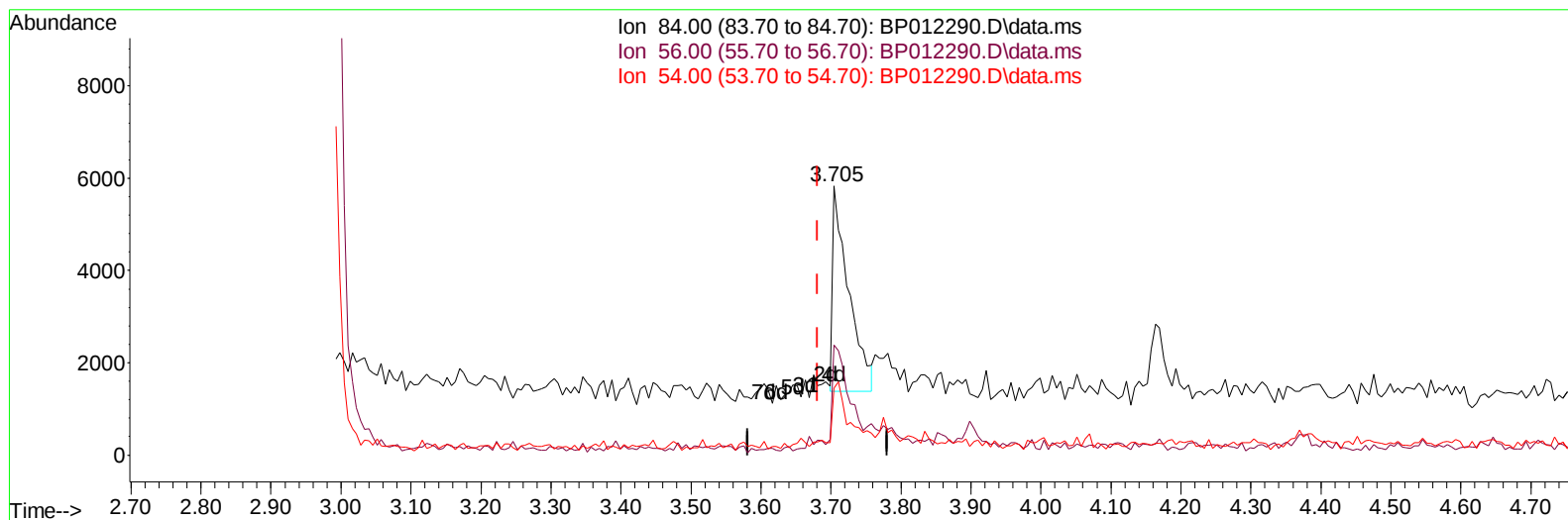
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(4) Pyridine-d5 (S)

3.705min (+ 0.023) 0.30 ng/ul m

response 7080

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	41.02#
54.00	29.70	24.84
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	327984	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1433465	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	791118	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.227	188	1371076	20.000	ng/ul	0.00
79) Chrysene-d12	21.321	240	1251910	20.000	ng/ul	0.00
88) Perylene-d12	23.715	264	1379918	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	15450	1.906	ng/uL	0.00
4) Pyridine-d5	3.705	84	7080m	0.303	ng/ul	0.02
7) Phenol-d5	6.987	99	721925	25.496	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.151	67	481586	28.045	ng/ul	0.00
11) 2-Chlorophenol-d4	7.346	132	568891	26.395	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	279467	12.699	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128	315917	29.774	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	319303	27.892	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.245	165	531430	24.748	ng/ul	0.00
31) 4-Chloroaniline-d4	10.769	131	610666	20.016	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	1665745	30.269	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1939351	30.874	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	151045	14.651	ng/ul	0.00
60) Fluorene-d10	15.463	176	1382831	29.208	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	81149	10.404	ng/ul	0.00
73) Anthracene-d10	17.327	188	1738394	28.926	ng/ul	0.00
81) Pyrene-d10	19.563	212	2041127	28.636	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	2114505	30.556	ng/ul	0.00
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	37161	1.254	ng/ul#	87

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

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