

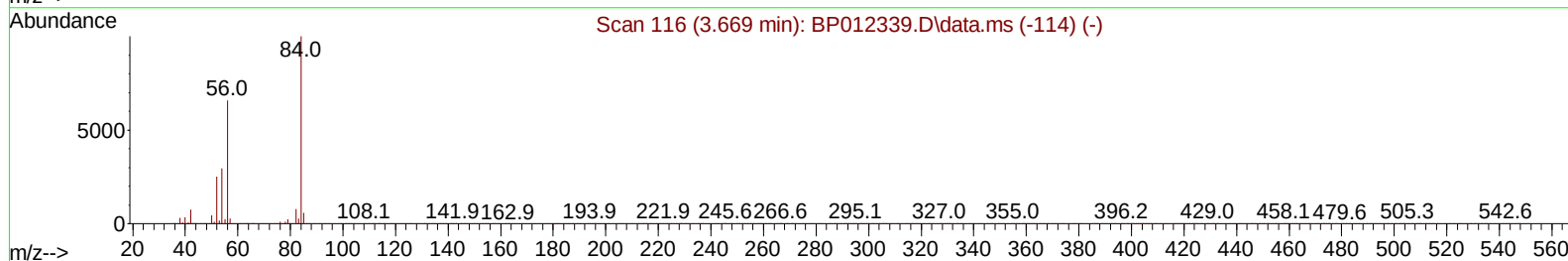
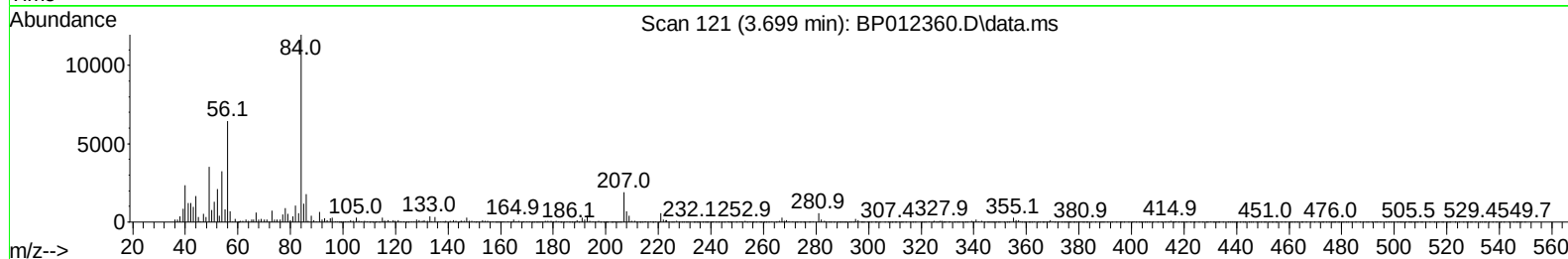
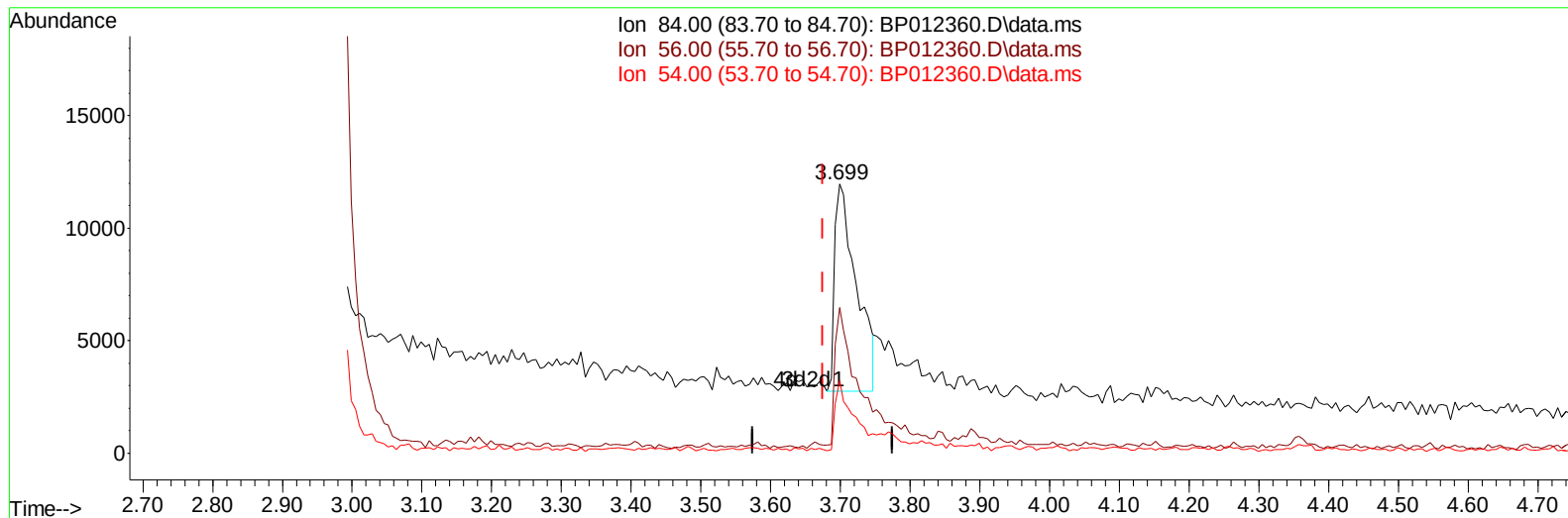
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012360.D
Acq On : 28 Oct 2022 01:13
Operator : CG/JU
Sample : N5144-05
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0BQ3

Manual IntegrationsAPPROVED

Quant Time: Oct 28 01:47:08 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 26 23:40:20 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/28/2022
Supervised By :mohammad ahmed 10/31/2022



TIC: BP012360.D\data.ms

(4) Pyridine-d5 (S)

3.699min (+ 0.024) 1.02 ng/u1

response 19771

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	54.04
54.00	29.80	27.09
0.00	0.00	0.00

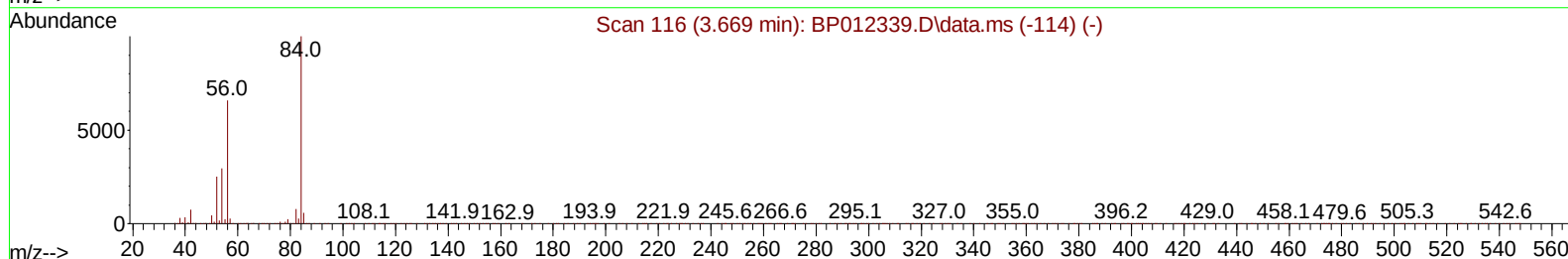
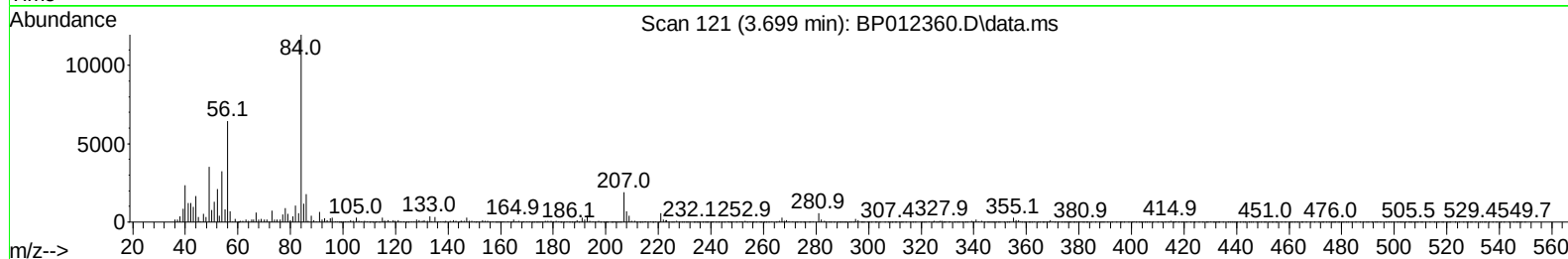
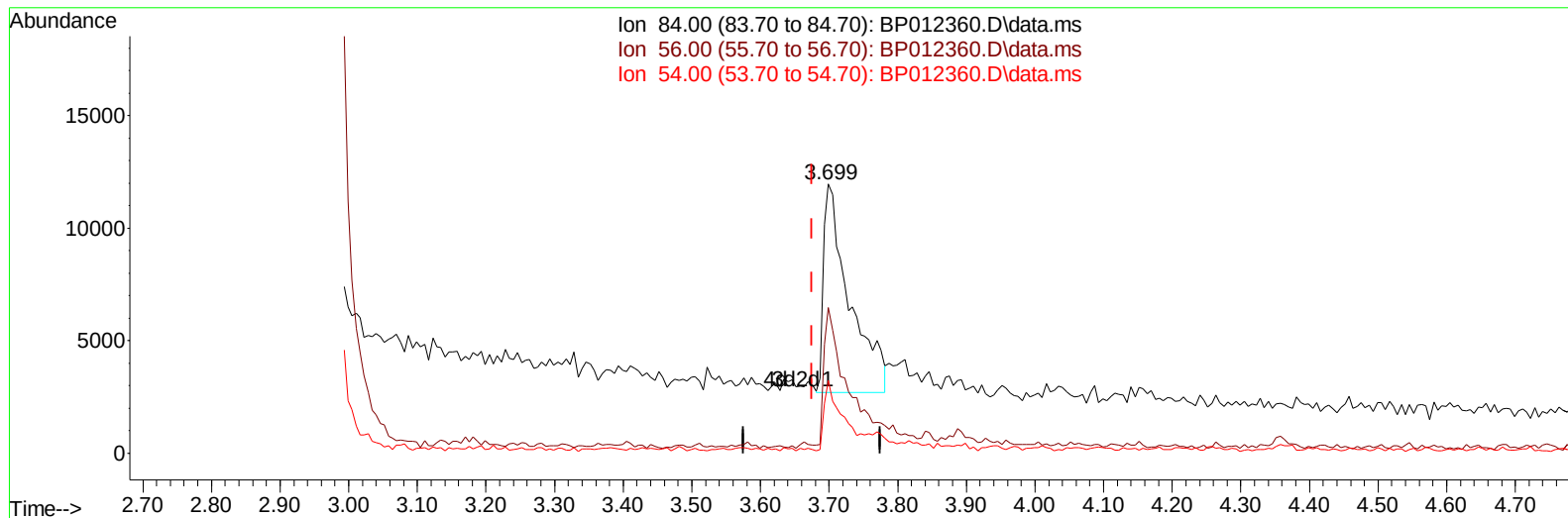
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(4) Pyridine-d5 (S)

3.699min (+ 0.024) 1.26 ng/ul m

response 24341

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	54.04
54.00	29.80	27.09
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	7.805	152	271355	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.610	136	1252360	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	841534	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	1842341	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	2046530	20.000	ng/ul	0.00
88) Perylene-d12	23.704	264	2012236	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	27359	4.060	ng/uL	0.00
4) Pyridine-d5	3.699	84	24341m	1.257	ng/ul	0.02
7) Phenol-d5	6.975	99	580145	23.837	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.140	67	359017	24.710	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.334	132	460517	25.910	ng/ul	-0.01
15) 4-Methylphenol-d8	8.522	113	301126	15.643	ng/ul	-0.01
21) Nitrobenzene-d5	8.975	128	241431	29.728	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.693	143	268041	31.799	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.234	165	453740	24.838	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	607265	22.326	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	1659752	28.811	ng/ul	-0.01
49) Acenaphthylene-d8	14.151	160	1849954	28.101	ng/ul	0.00
54) 4-Nitrophenol-d4	14.675	143	267001	24.489	ng/ul	0.00
60) Fluorene-d10	15.451	176	1464307	29.689	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.586	200	192806	20.101	ng/ul	0.00
73) Anthracene-d10	17.316	188	2322115	29.252	ng/ul	0.00
81) Pyrene-d10	19.557	212	3001771	29.214	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	3076324	31.289	ng/ul	0.00

Target Compounds Qvalue

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

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