

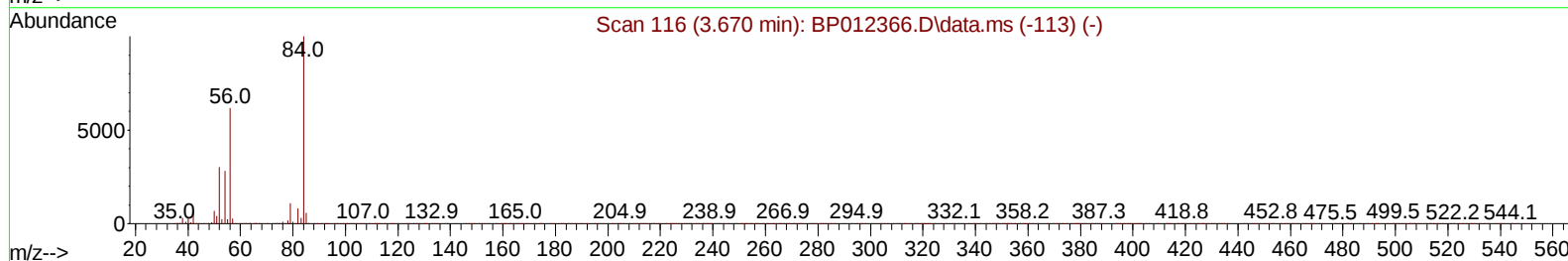
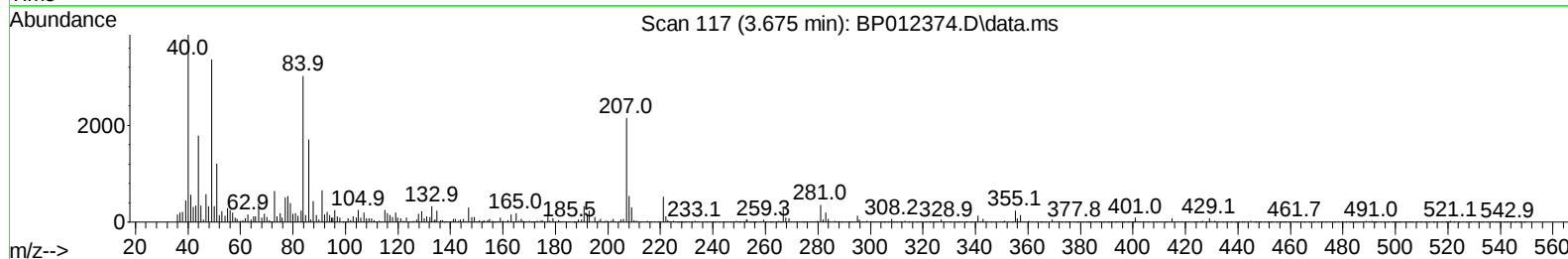
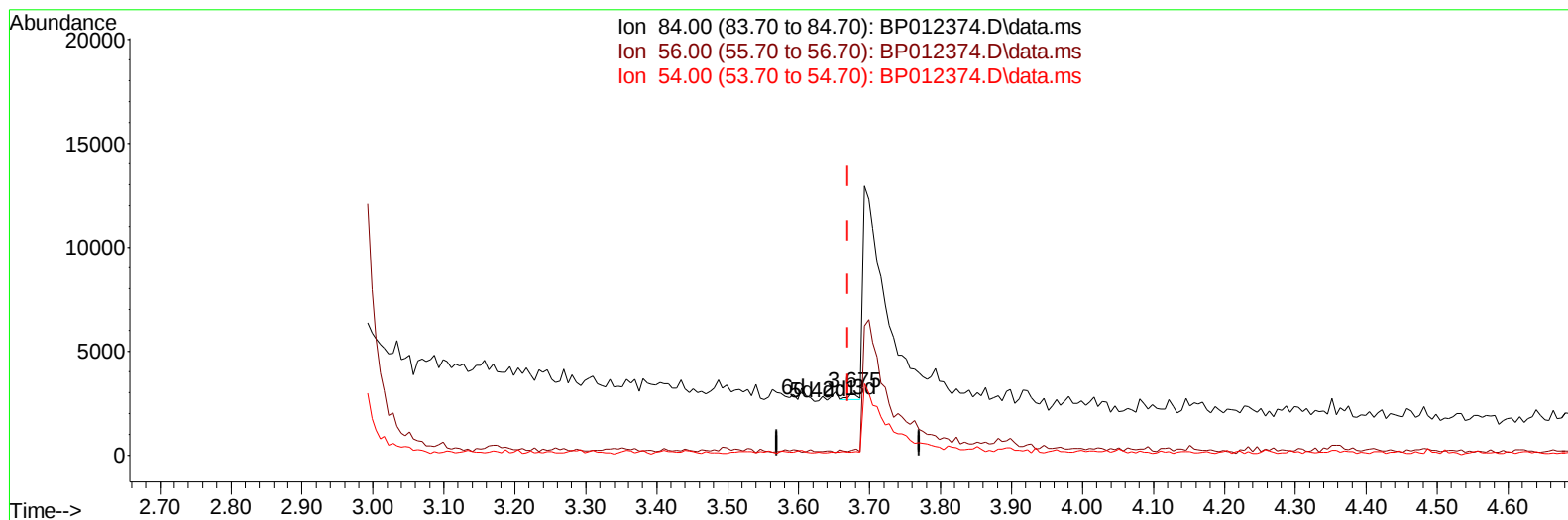
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012374.D
 Acq On : 28 Oct 2022 12:57
 Operator : CG/JU
 Sample : N5232-14
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA82

Manual IntegrationsAPPROVED

Quant Time: Oct 28 23:21:05 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Oct 28 06:05:59 2022
 Response via : Initial Calibration

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



TIC: BP012374.D\data.ms

(4) Pyridine-d5 (S)

3.675min (+ 0.005) 0.02 ng/u1

response	294	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	8.18#
54.00	29.80	4.24#
0.00	0.00	0.00

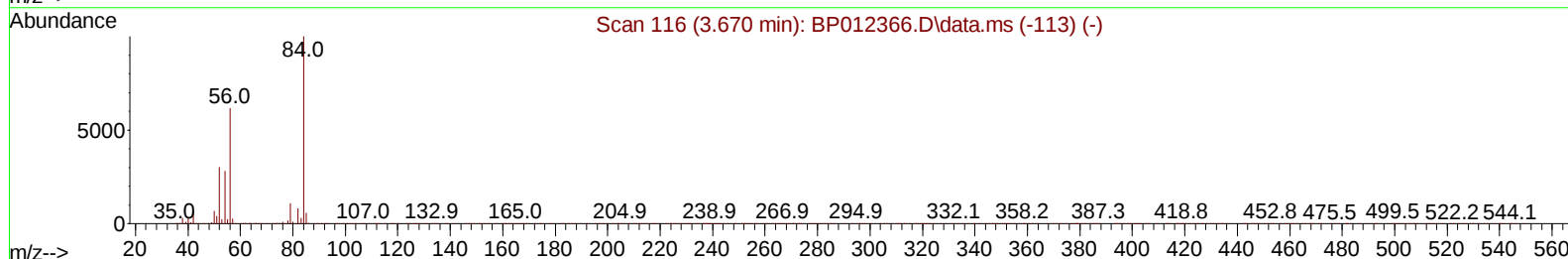
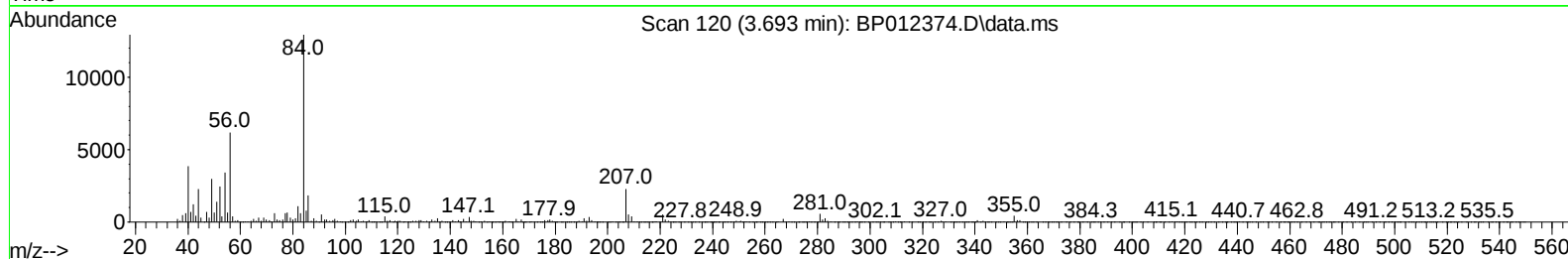
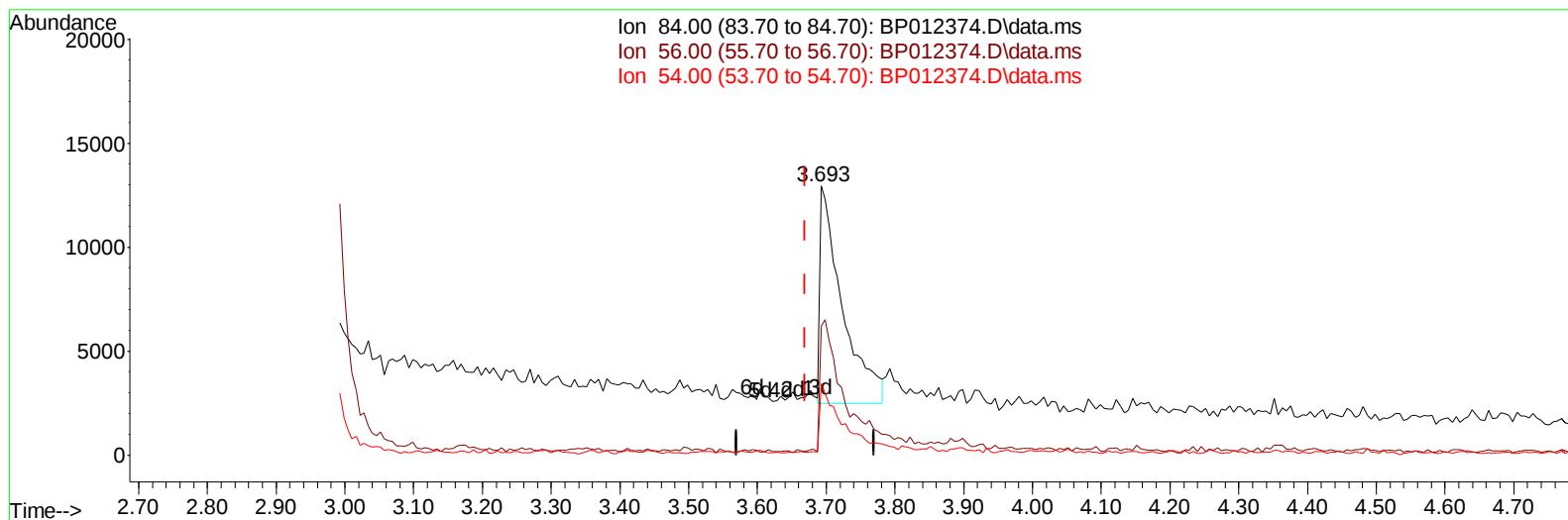
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(4) Pyridine-d5 (S)

3.693min (+ 0.023) 1.49 ng/u1 m

response 23672

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	47.96#
54.00	29.80	26.55
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	7.804	152	222902	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	1022965	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	685788	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.215	188	1553822	20.000	ng/ul	0.00
79) Chrysene-d12	21.309	240	1848432	20.000	ng/ul	0.00
88) Perylene-d12	23.703	264	1802824	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	25314	4.573	ng/uL	0.00
4) Pyridine-d5	3.693	84	23672m	1.489	ng/ul	0.02
7) Phenol-d5	6.981	99	160014	8.004	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	432632	36.249	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	440949	30.202	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	301401	19.061	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	298312	44.969	ng/ul	0.00
24) 2-Nitrophenol-d4	9.692	143	308607	44.821	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	516928	34.643	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	69795	3.141	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	1896821	40.404	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	2177763	40.593	ng/ul	0.00
54) 4-Nitrophenol-d4	14.686	143	56745	6.387	ng/ul	0.01
60) Fluorene-d10	15.451	176	1698413	42.255	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	267900	33.116	ng/ul	0.00
73) Anthracene-d10	17.315	188	2894050	43.227	ng/ul	0.00
81) Pyrene-d10	19.557	212	3646509	39.293	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.550	264	3815937	43.319	ng/ul	0.00

Target Compounds Qvalue

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

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