

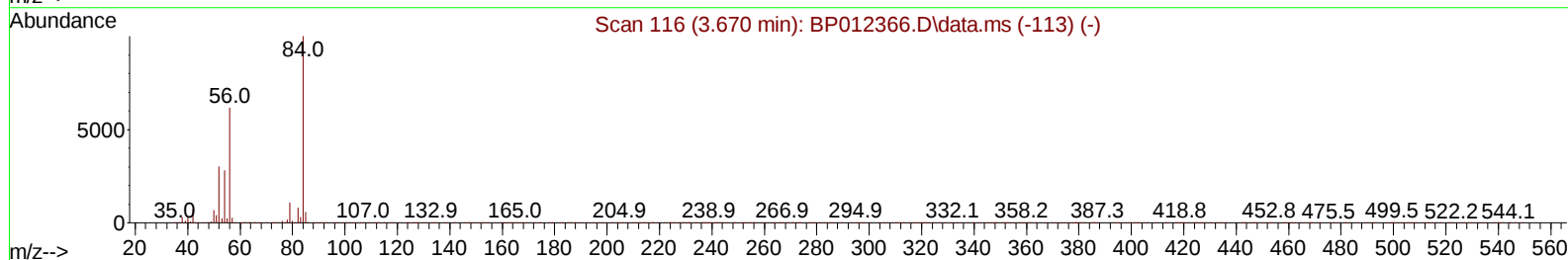
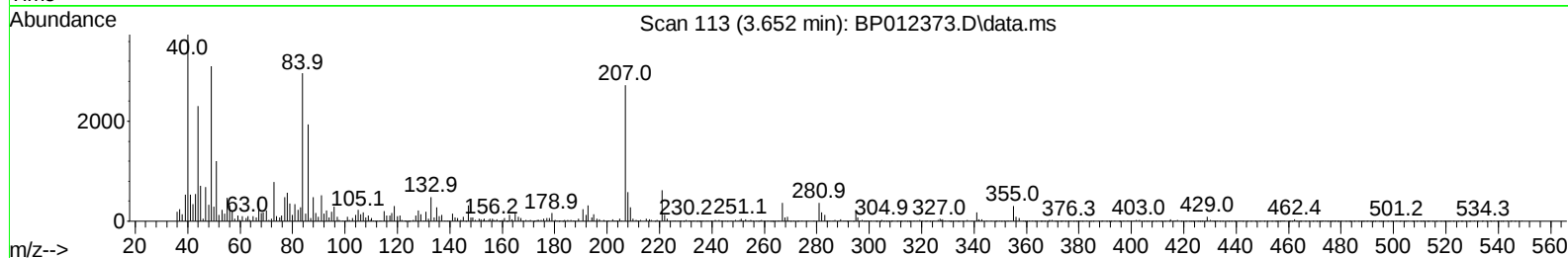
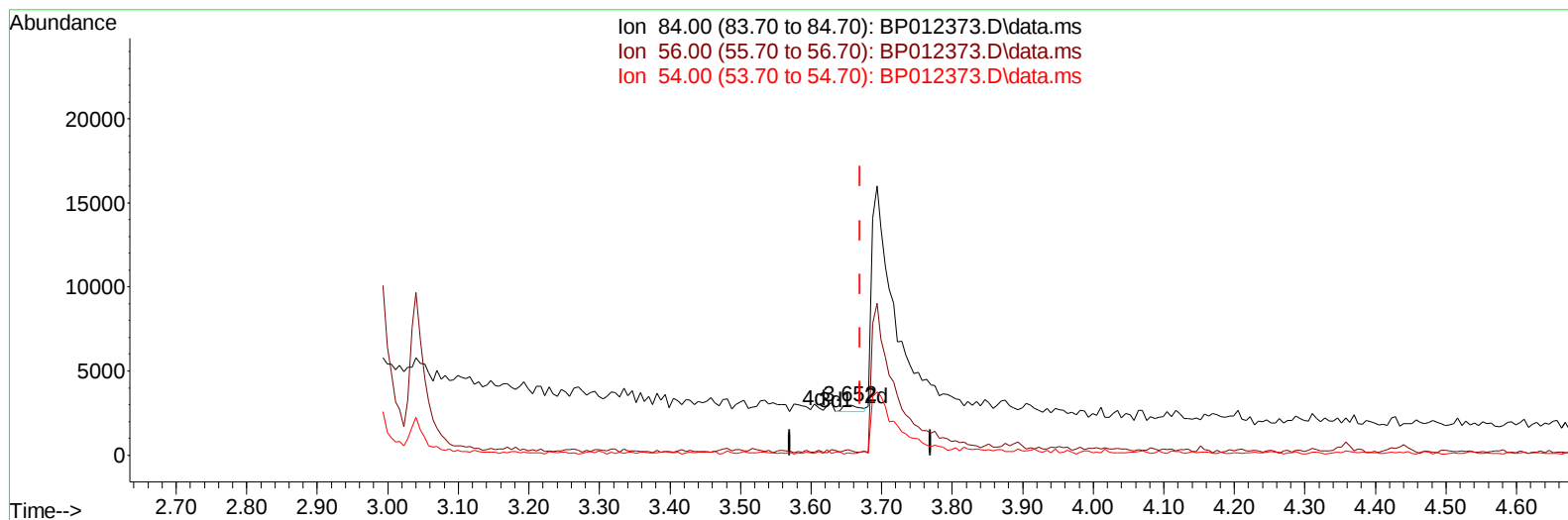
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012373.D
 Acq On : 28 Oct 2022 12:23
 Operator : CG/JU
 Sample : N5232-13
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA71

Manual IntegrationsAPPROVED

Quant Time: Oct 28 23:20:56 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: BP012373.D\data.ms

(4) Pyridine-d5 (S)

3.652min (-0.018) 0.04 ng/u1

response 661

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	10.67#
54.00	29.80	5.10#
0.00	0.00	0.00

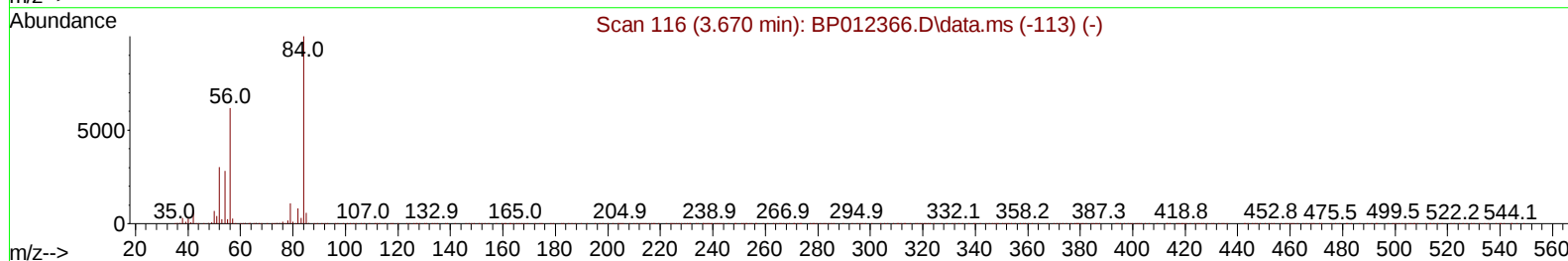
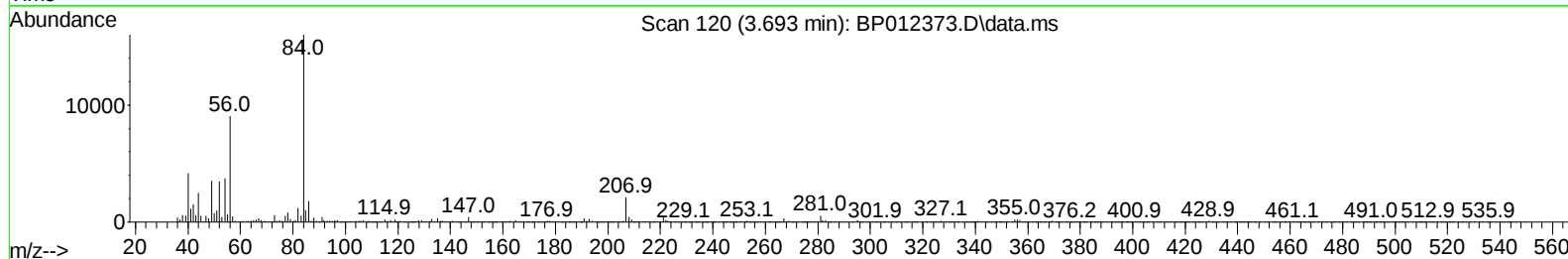
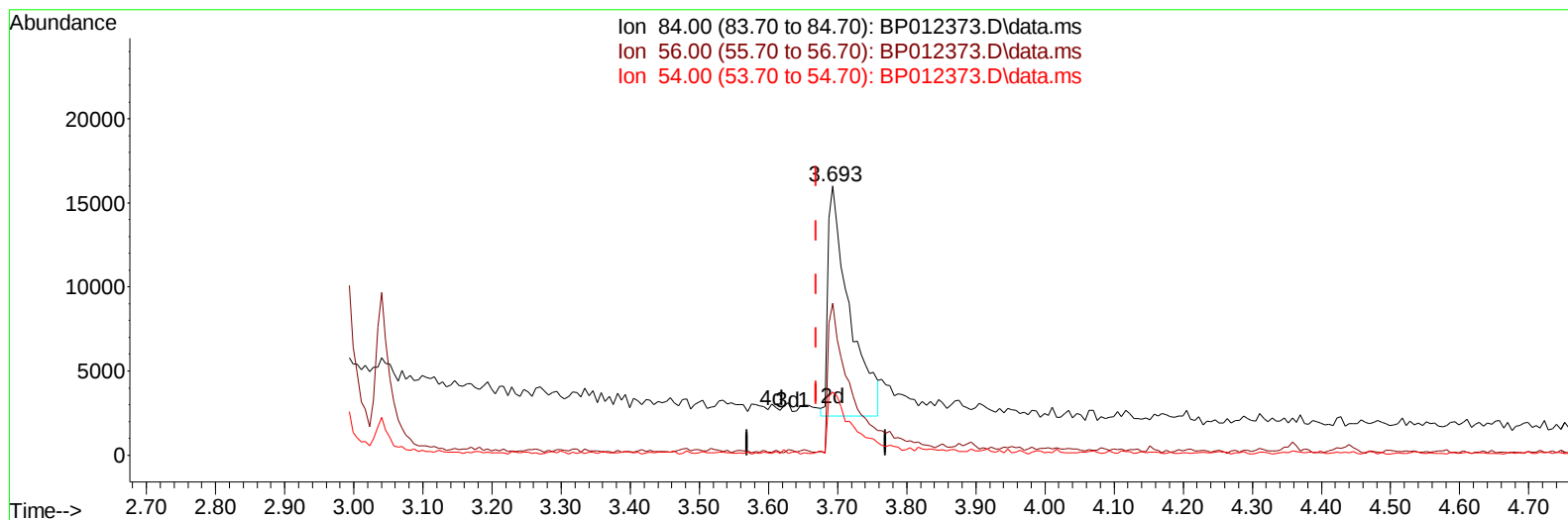
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(4) Pyridine-d5 (S)

3.693min (+ 0.024) 1.62 ng/ul m

response 29415

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	56.48
54.00	29.80	23.38#
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	253892	20.000	ng/ul	0.00
20) Naphthalene-d8	10.605	136	1161461	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	751975	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	1697664	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1965437	20.000	ng/ul	0.00
88) Perylene-d12	23.704	264	1860639	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	25414	4.030	ng/uL	0.00
4) Pyridine-d5	3.693	84	29415m	1.624	ng/ul	0.02
7) Phenol-d5	6.981	99	155766	6.840	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	410532	30.199	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	420938	25.312	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	302157	16.777	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	280209	37.203	ng/ul	0.00
24) 2-Nitrophenol-d4	9.693	143	290082	37.107	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	526546	31.079	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	59237	2.348	ng/ul	0.01
46) Dimethylphthalate-d6	13.869	166	1961609	38.107	ng/ul	0.00
49) Acenaphthylene-d8	14.146	160	2113248	35.923	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	69270	7.110	ng/ul	0.00
60) Fluorene-d10	15.451	176	1718295	38.987	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.587	200	292478	33.090	ng/ul	0.00
73) Anthracene-d10	17.316	188	2951911	40.355	ng/ul	0.00
81) Pyrene-d10	19.557	212	3706740	37.564	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	3743282	41.174	ng/ul	0.00

Target Compounds Qvalue

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

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