

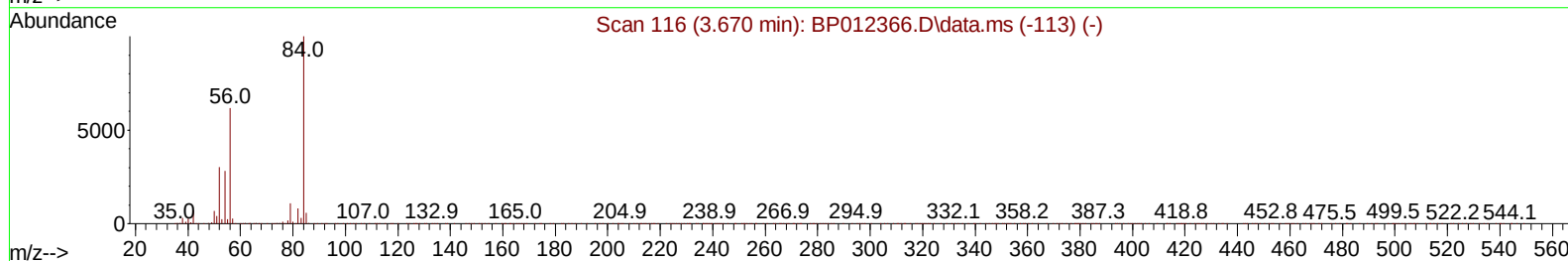
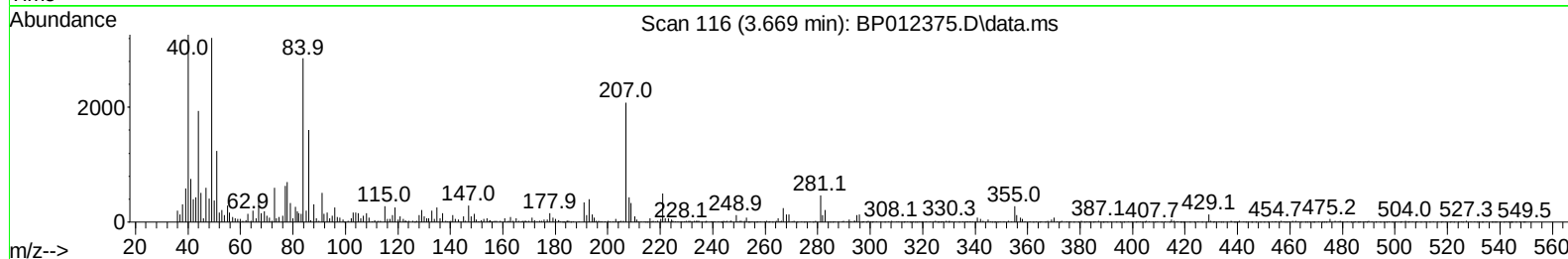
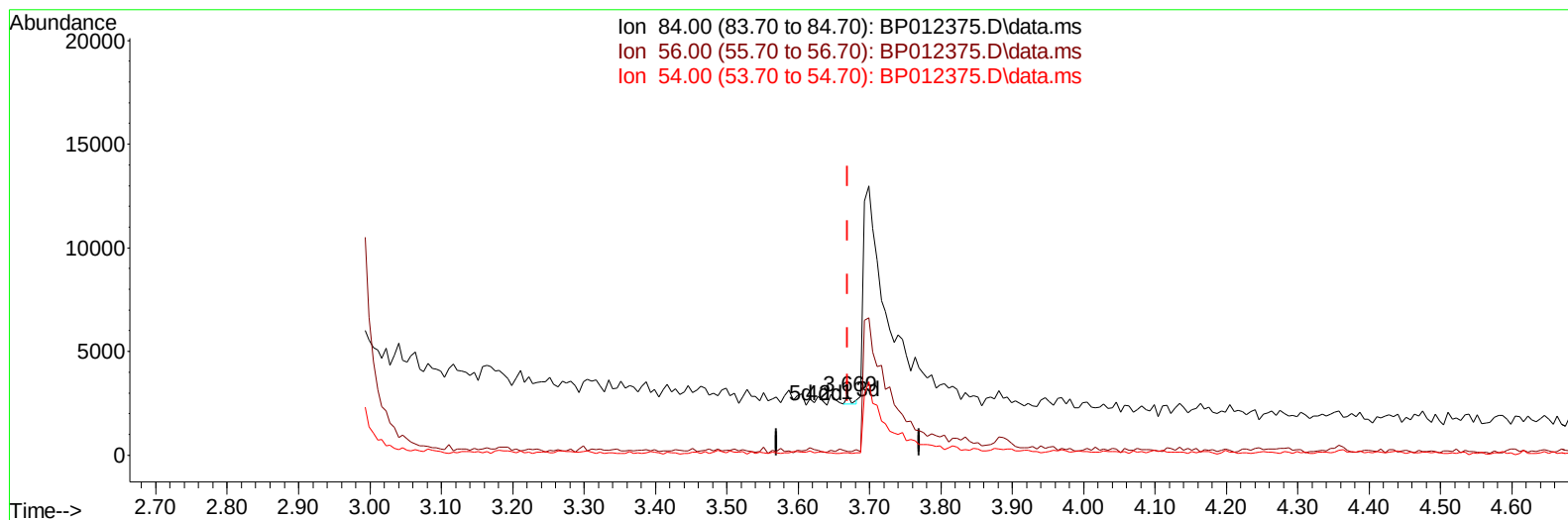
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
 Data File : BP012375.D
 Acq On : 28 Oct 2022 13:31
 Operator : CG/JU
 Sample : N5232-15
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHA90

Manual IntegrationsAPPROVED

Quant Time: Oct 28 23:21:15 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: BP012375.D\data.ms

(4) Pyridine-d5 (S)

3.669min (-0.000) 0.01 ng/uI

response	215	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	5.90#
54.00	29.80	4.28#
0.00	0.00	0.00

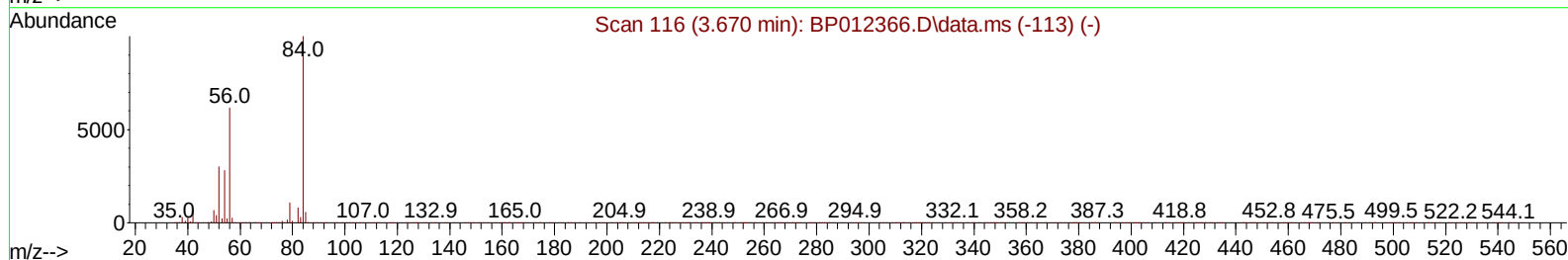
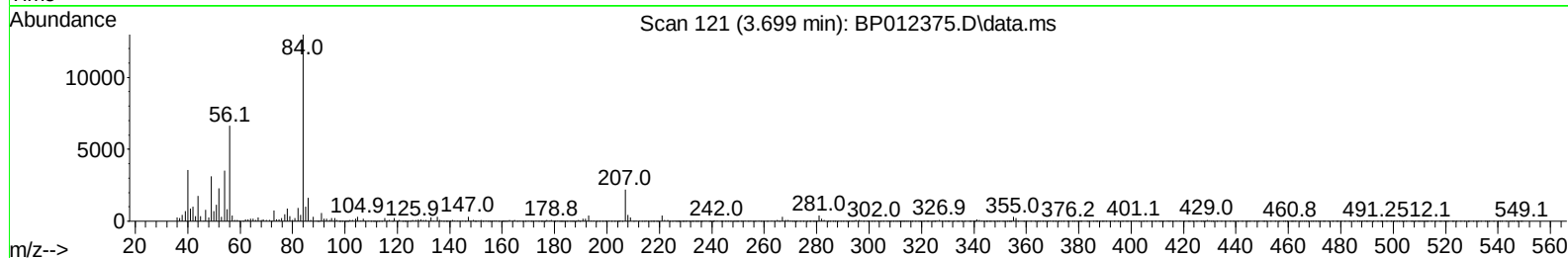
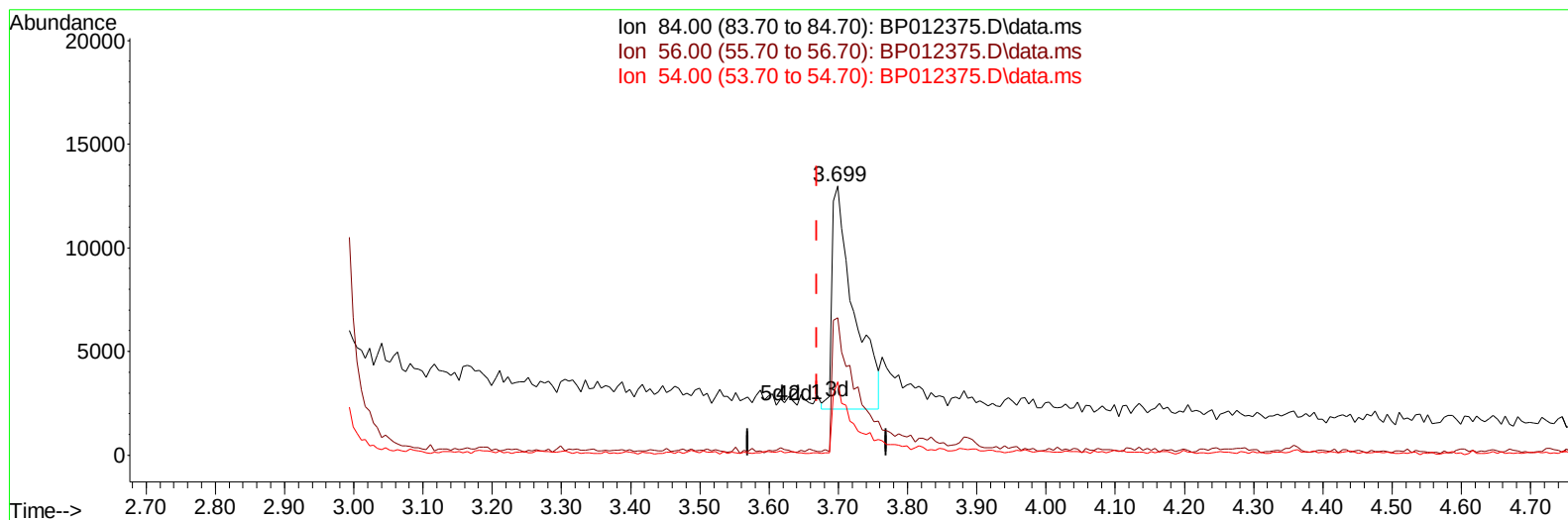
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(4) Pyridine-d5 (S)

3.699min (+ 0.029) 1.47 ng/ul m

response 23384

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	51.14#
54.00	29.80	27.26
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	223634	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	1012039	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	672282	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.216	188	1547438	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1827803	20.000	ng/ul	0.00
88) Perylene-d12	23.704	264	1777034	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	24012	4.323	ng/uL	0.00
4) Pyridine-d5	3.699	84	23384m	1.466	ng/ul	0.03
7) Phenol-d5	6.981	99	151917	7.574	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	407786	34.056	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	414742	28.314	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	283533	17.872	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	275627	41.998	ng/ul	0.00
24) 2-Nitrophenol-d4	9.693	143	279068	40.968	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	478212	32.394	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	49273	2.242	ng/ul	0.01
46) Dimethylphthalate-d6	13.875	166	1790552	38.907	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	2076053	39.474	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	53151	6.102	ng/ul	0.01
60) Fluorene-d10	15.451	176	1603756	40.702	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.586	200	249463	30.964	ng/ul	0.00
73) Anthracene-d10	17.316	188	2761806	41.422	ng/ul	0.00
81) Pyrene-d10	19.557	212	3466864	37.779	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.551	264	3628948	41.795	ng/ul	0.00

Target Compounds Qvalue

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

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