

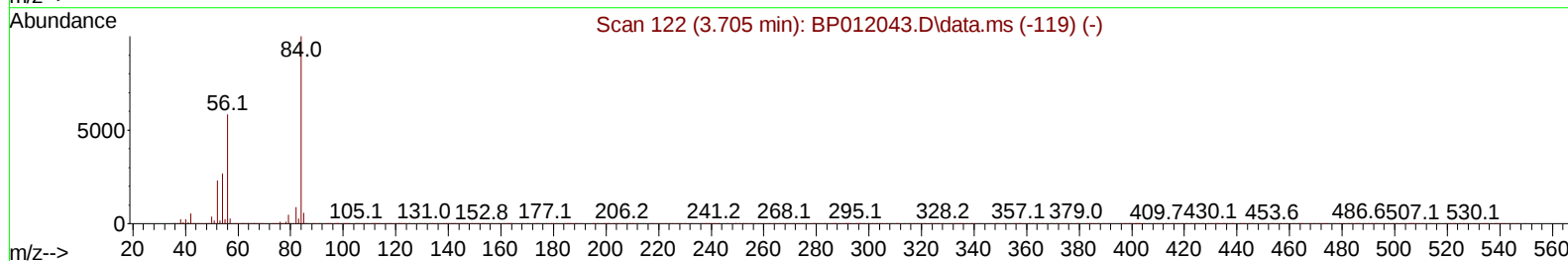
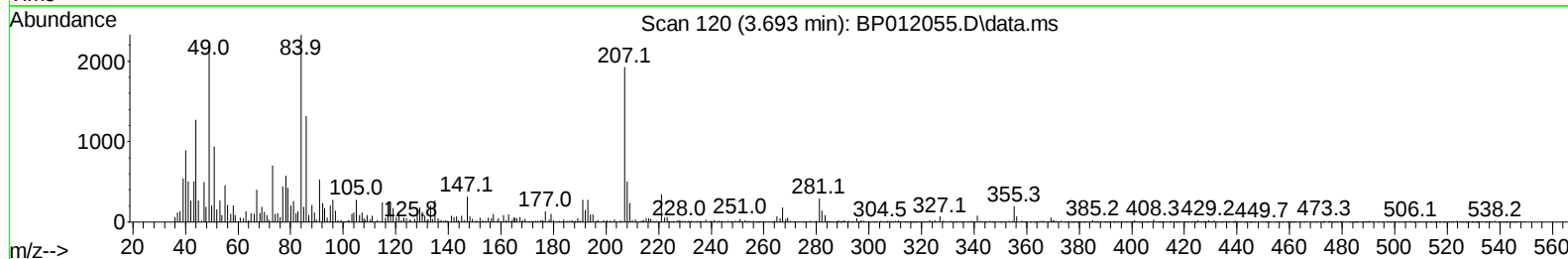
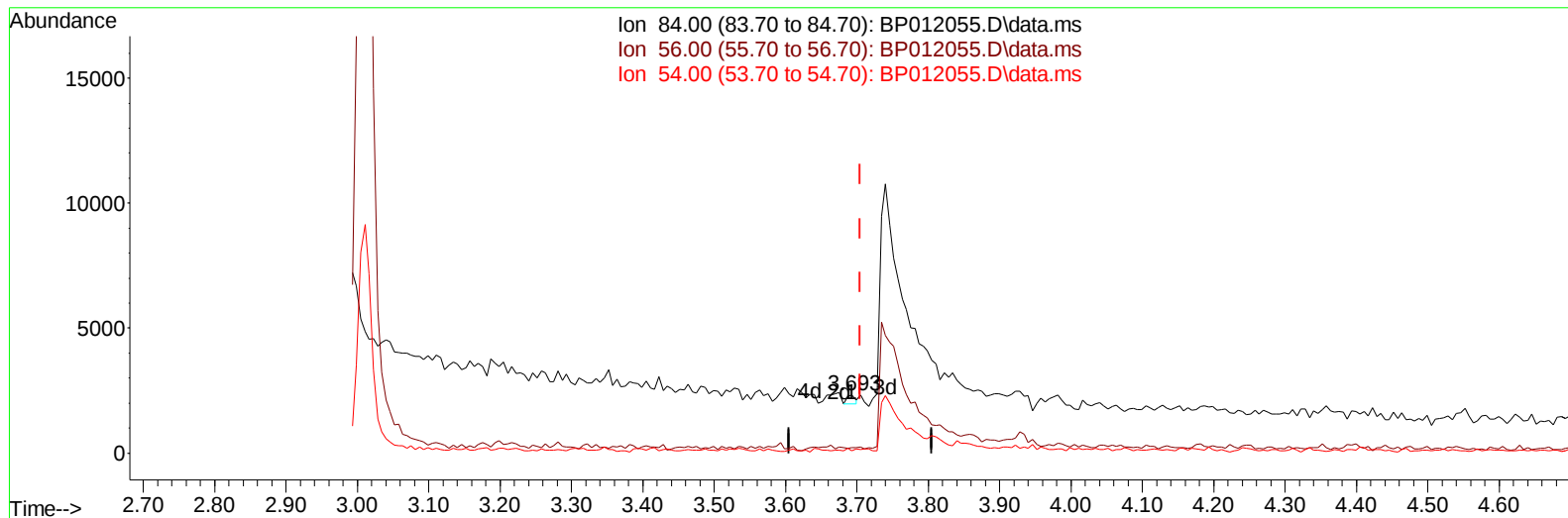
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101122\
 Data File : BP012055.D
 Acq On : 11 Oct 2022 23:43
 Operator : CG/JU
 Sample : N4991-02
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BE9

Manual IntegrationsAPPROVED

Quant Time: Oct 12 00:25:25 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Oct 11 23:54:43 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/12/2022
 Supervised By :mohammad ahmed 10/14/2022



TIC: BP012055.D\data.ms

(4) Pyridine-d5 (S)

3.693min (-0.012) 0.02 ng/u1

response 276

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	9.15#
54.00	26.00	3.65#
0.00	0.00	0.00

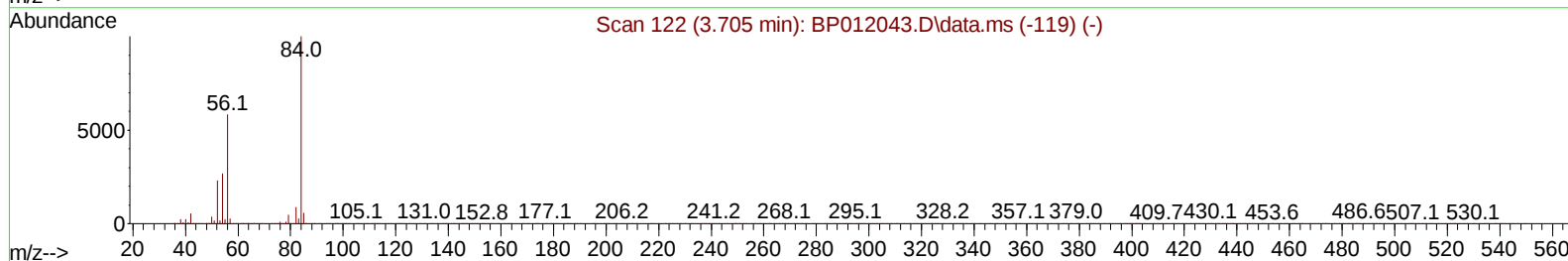
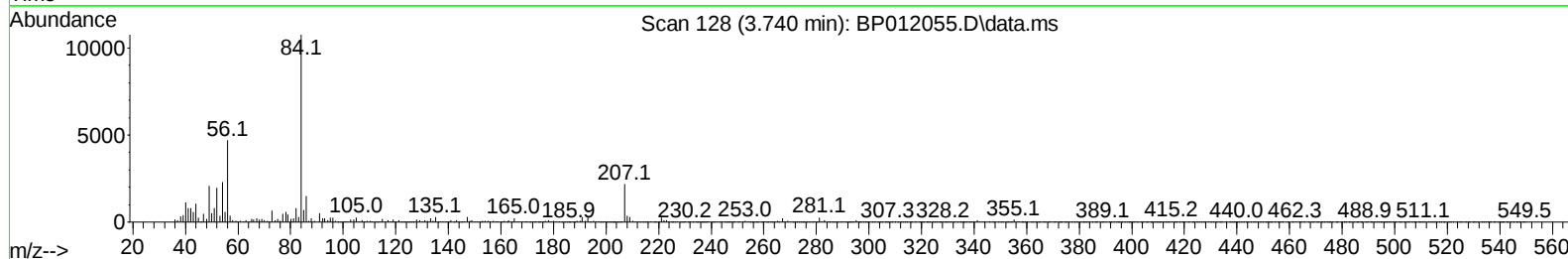
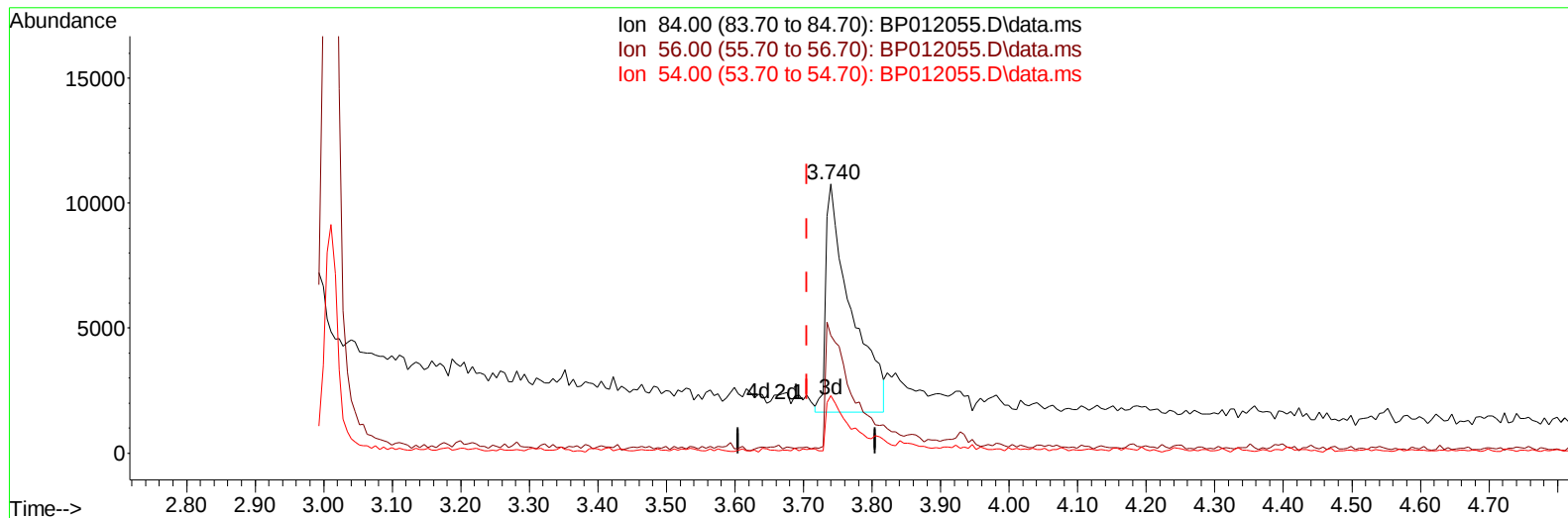
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(4) Pyridine-d5 (S)

3.740min (+ 0.035) 1.42 ng/ul m

response 23128

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	43.67#
54.00	26.00	21.25
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	7.858	152	240803	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	956879	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.504	164	546994	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.263	188	1165584	20.000	ng/ul	0.00
79) Chrysene-d12	21.351	240	1302929	20.000	ng/ul	0.00
88) Perylene-d12	23.769	264	1408127	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	18498	3.029	ng/uL	0.00
4) Pyridine-d5	3.740	84	23128m	1.424	ng/ul	0.04
7) Phenol-d5	7.022	99	233833	12.074	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	179556	16.334	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	247020	16.069	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	149530	10.052	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	116845	17.290	ng/ul	0.00
24) 2-Nitrophenol-d4	9.746	143	119752	17.476	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	221320	16.665	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	283618	14.492	ng/ul	0.00
46) Dimethylphthalate-d6	13.916	166	705144	20.355	ng/ul	0.00
49) Acenaphthylene-d8	14.198	160	794430	18.545	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	65394	9.615	ng/ul	0.00
60) Fluorene-d10	15.498	176	639559	20.406	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.622	200	62838	10.210	ng/ul	0.00
73) Anthracene-d10	17.363	188	1056091	20.866	ng/ul	0.00
81) Pyrene-d10	19.598	212	1421416	21.295	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.610	264	1498139	21.886	ng/ul	0.00

Target Compounds Qvalue

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

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