

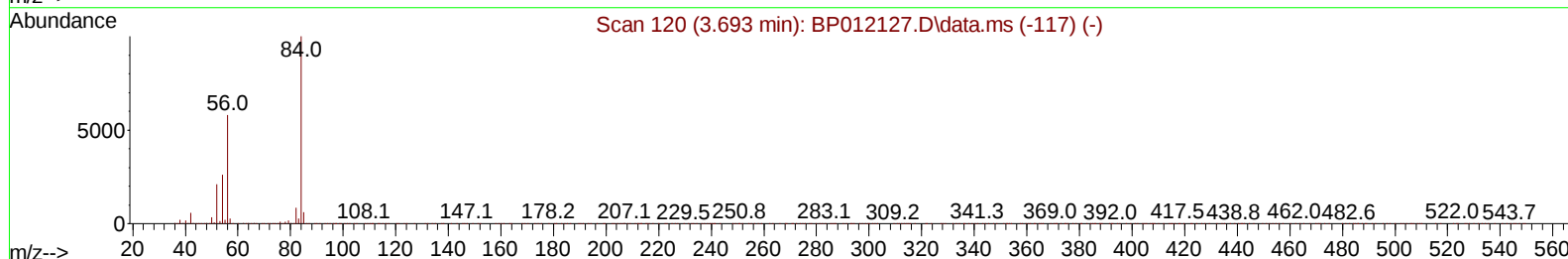
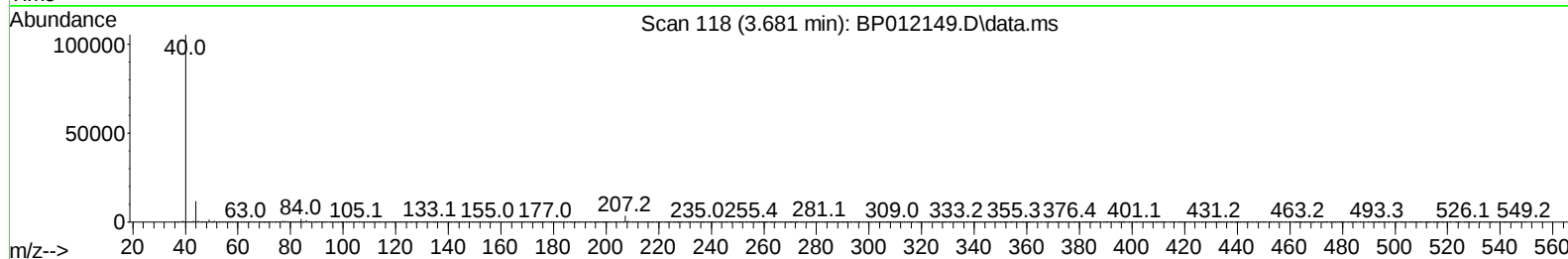
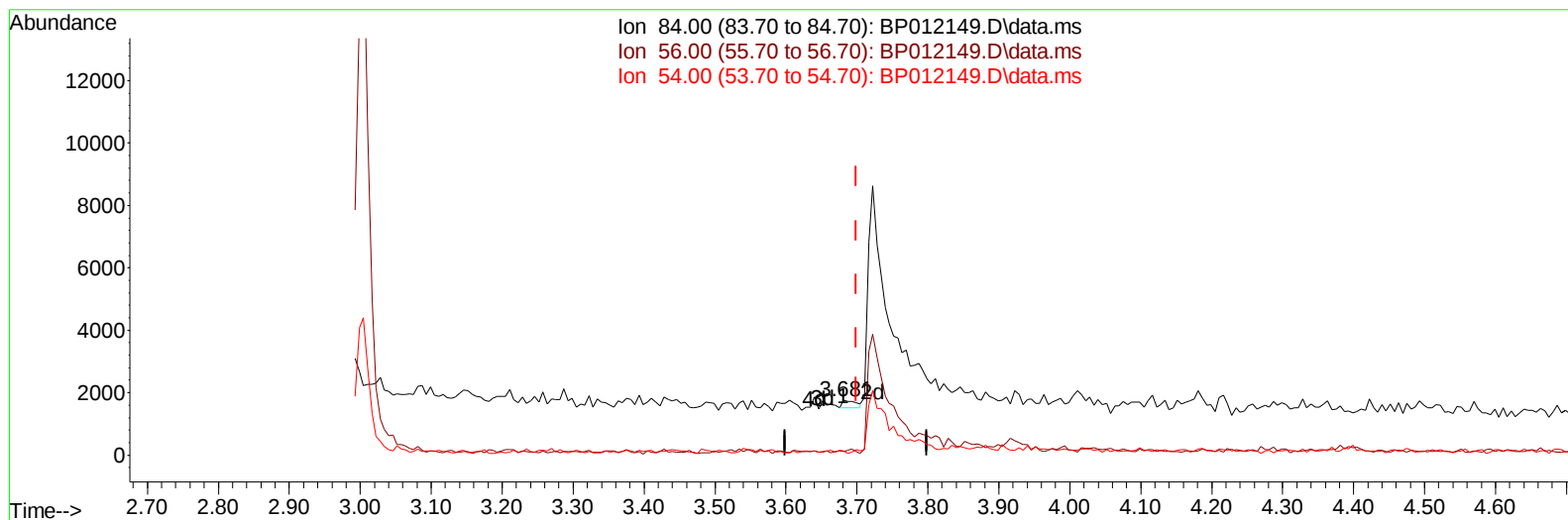
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101422\
 Data File : BP012149.D
 Acq On : 14 Oct 2022 22:10
 Operator : CG/JU
 Sample : N4984-01
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BGNC3

Manual IntegrationsAPPROVED

Quant Time: Oct 14 23:07:43 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 12 23:03:21 2022
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/15/2022
 Supervised By :mohammad ahmed 10/16/2022



TIC: BP012149.D\data.ms

(4) Pyridine-d5 (S)

3.681min (-0.018) 0.02 ng/u1

response 324

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	6.96#
54.00	26.00	7.64#
0.00	0.00	0.00

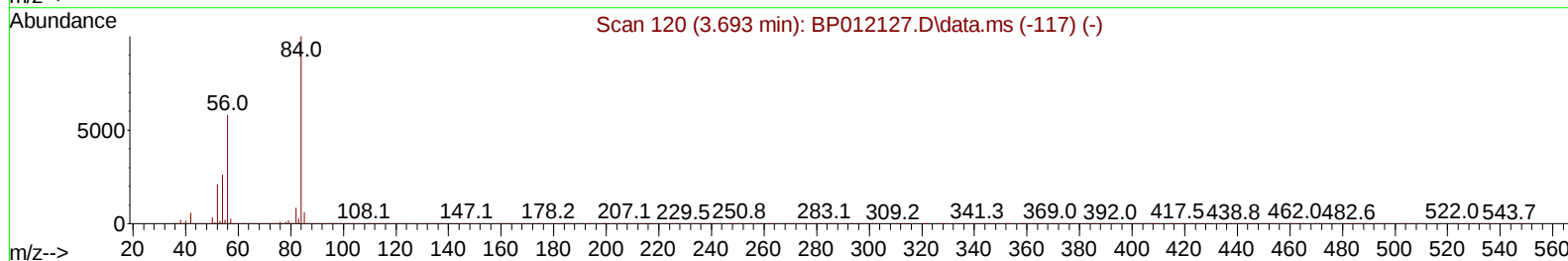
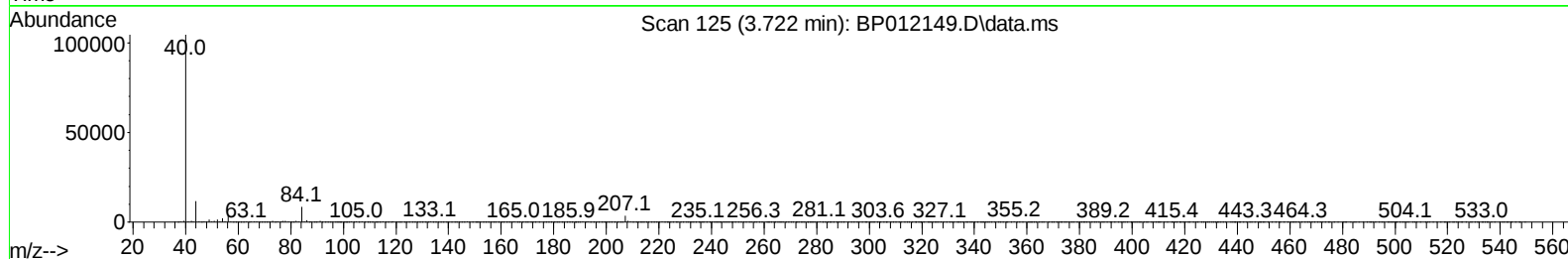
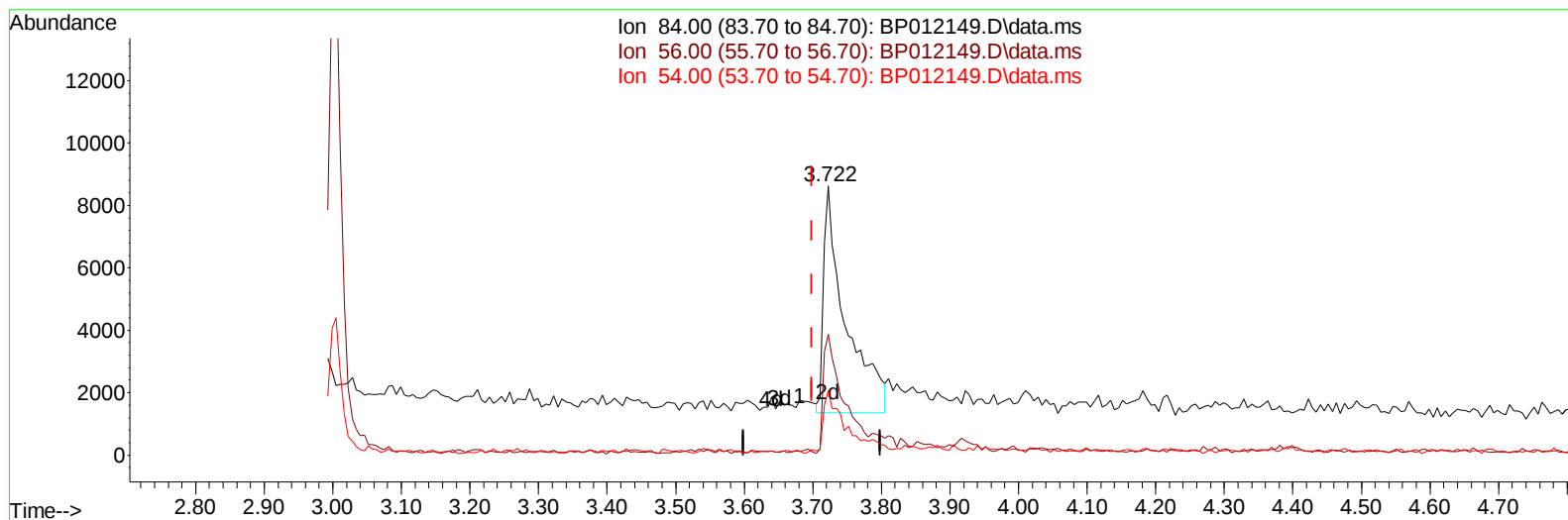
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TIC: BP012149.D\data.ms

(4) Pyridine-d5 (S)

3.722min (+ 0.023) 0.99 ng/u1 m

response 16214

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	57.80	45.01#
54.00	26.00	23.90
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	7.840	152	243633	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.645	136	992311	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.492	164	572614	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	1223381	20.000	ng/ul	0.00
79) Chrysene-d12	21.345	240	1252699	20.000	ng/ul	0.00
88) Perylene-d12	23.757	264	1309180	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	10540	1.706	ng/uL	0.00
4) Pyridine-d5	3.722	84	16214m	0.987	ng/ul	0.02
7) Phenol-d5	7.010	99	318743	16.267	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	194637	17.500	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.375	132	260426	16.744	ng/ul	0.00
15) 4-Methylphenol-d8	8.557	113	242779	16.131	ng/ul	0.00
21) Nitrobenzene-d5	9.010	128	120211	17.153	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.734	143	119549	16.823	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.269	165	210058	15.252	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.793	131	296211	14.595	ng/ul	-0.01
46) Dimethylphthalate-d6	13.904	166	638905	17.618	ng/ul	-0.01
49) Acenaphthylene-d8	14.186	160	759859	16.945	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	62188	8.735	ng/ul	0.00
60) Fluorene-d10	15.486	176	612265	18.661	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.616	200	40043	6.199	ng/ul	0.00
73) Anthracene-d10	17.351	188	1003469	18.889	ng/ul	0.00
81) Pyrene-d10	19.586	212	1186978	18.495	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.598	264	1151516	18.094	ng/ul	-0.01

Target Compounds Qvalue

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

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