

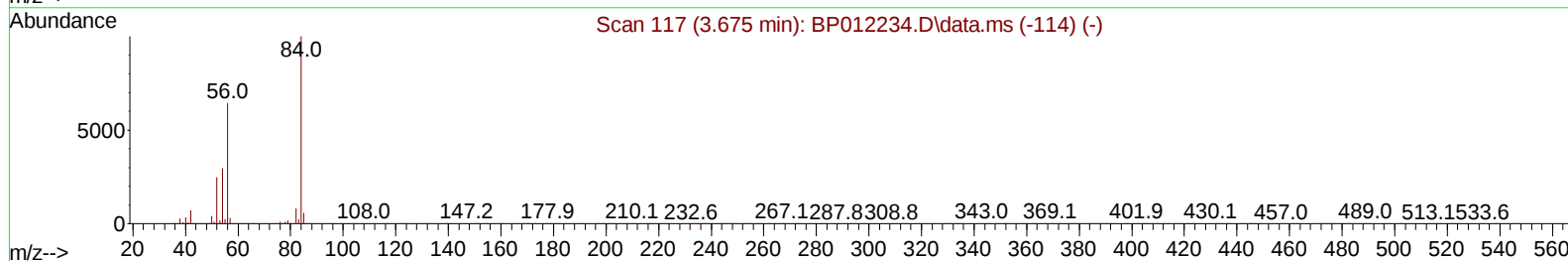
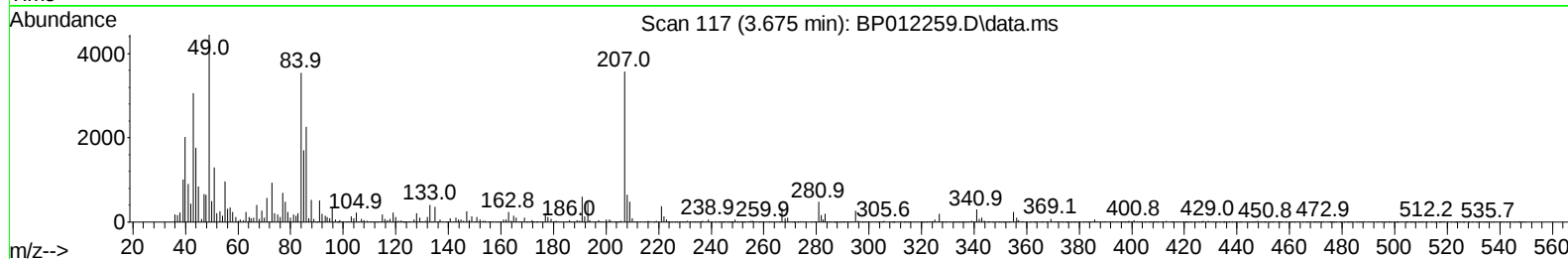
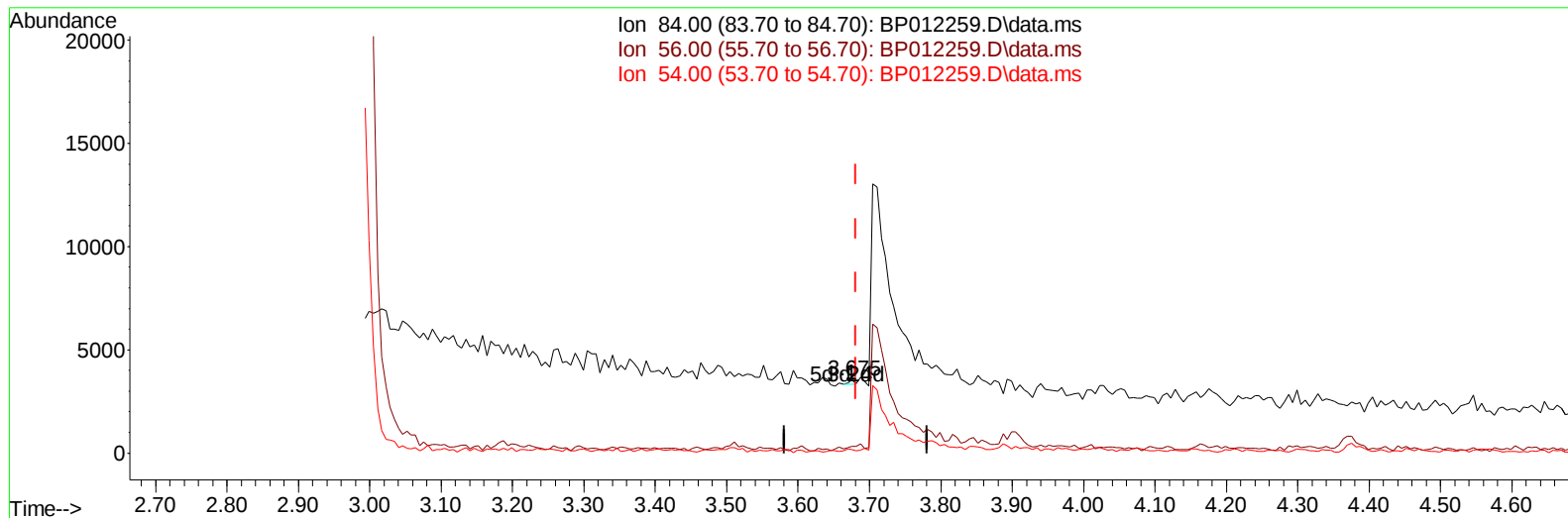
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012259.D
 Acq On : 21 Oct 2022 22:58
 Operator : CG/JU
 Sample : N5064-05
 Misc :
 ALS Vial : 54 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BL3

Manual IntegrationsAPPROVED

Quant Time: Oct 22 01:13:28 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 20 00:51:41 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/24/2022
 Supervised By :mohammad ahmed 10/24/2022



TIC: BP012259.D\data.ms

(4) Pyridine-d5 (S)

3.675min (-0.006) 0.00 ng/u1

response 124

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	8.75#
54.00	29.70	4.46#
0.00	0.00	0.00

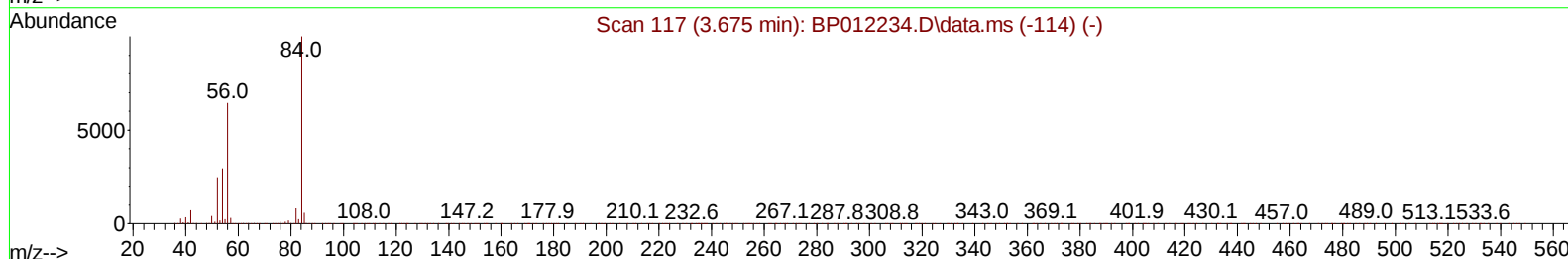
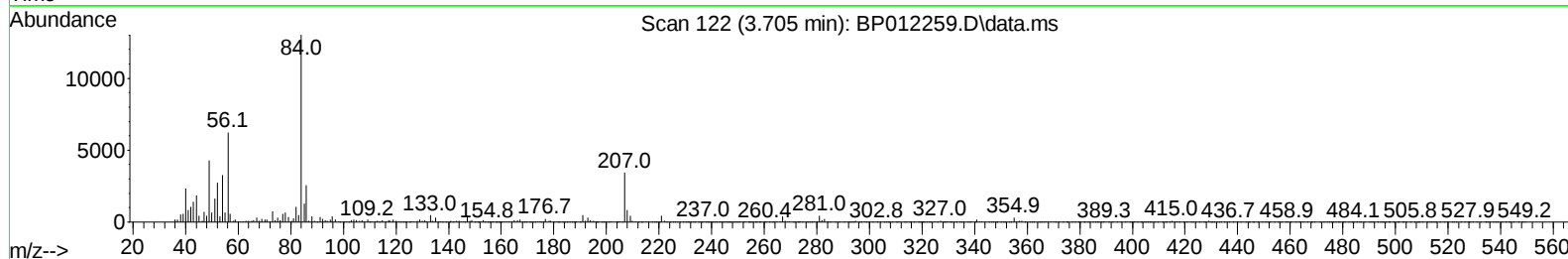
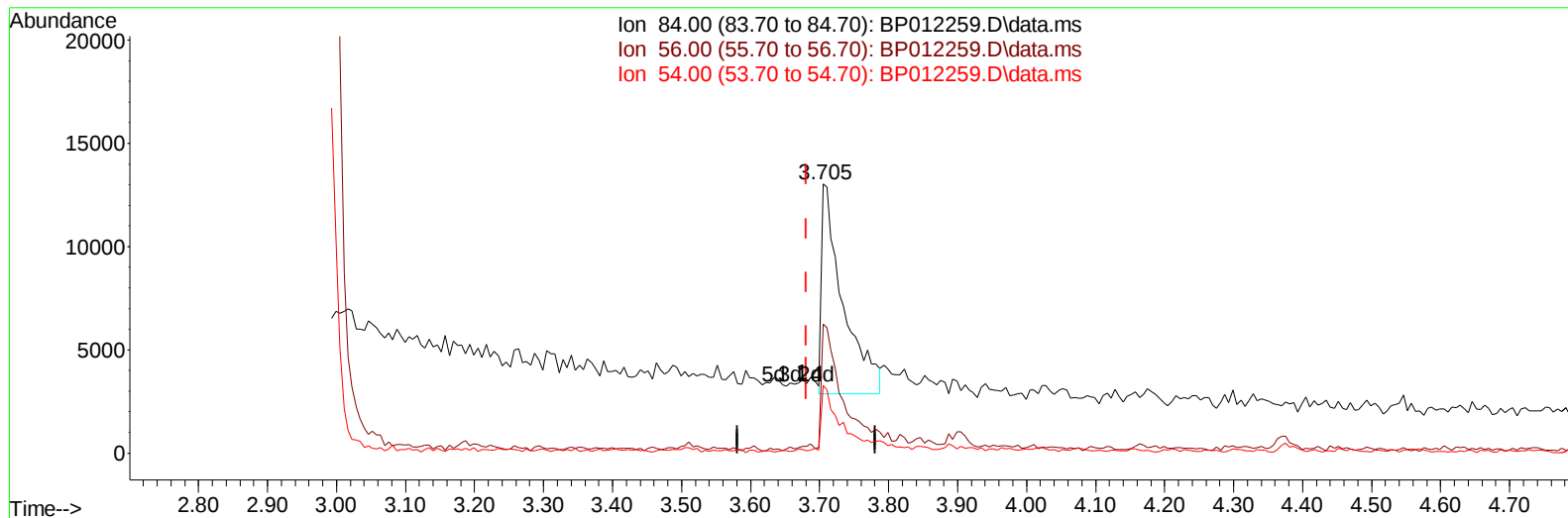
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(4) Pyridine-d5 (S)

3.705min (+ 0.024) 0.76 ng/ul m

response 22136

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	47.91#
54.00	29.70	25.05
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	7.822	152	406752	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1710677	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.469	164	979923	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.233	188	1710939	20.000	ng/ul	0.00
79) Chrysene-d12	21.322	240	1436951	20.000	ng/ul	-0.01
88) Perylene-d12	23.721	264	1561954	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	32192	3.203	ng/uL	0.00
4) Pyridine-d5	3.705	84	22136m	0.764	ng/ul	0.02
7) Phenol-d5	6.993	99	633674	18.045	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	397956	18.687	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	506403	18.946	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	341049	12.496	ng/ul	-0.01
21) Nitrobenzene-d5	8.987	128	257003	20.296	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.710	143	280825	20.556	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	476057	18.577	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.769	131	504633	13.860	ng/ul	-0.01
46) Dimethylphthalate-d6	13.887	166	1427220	20.938	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	1622436	20.852	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.692	143	160348	12.556	ng/ul	0.00
60) Fluorene-d10	15.469	176	1211180	20.654	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	92386	9.492	ng/ul	0.00
73) Anthracene-d10	17.328	188	1606764	21.425	ng/ul	-0.01
81) Pyrene-d10	19.569	212	1668188	20.390	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.568	264	1647890	21.038	ng/ul	-0.02
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	39043	1.063	ng/ul	94

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

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