

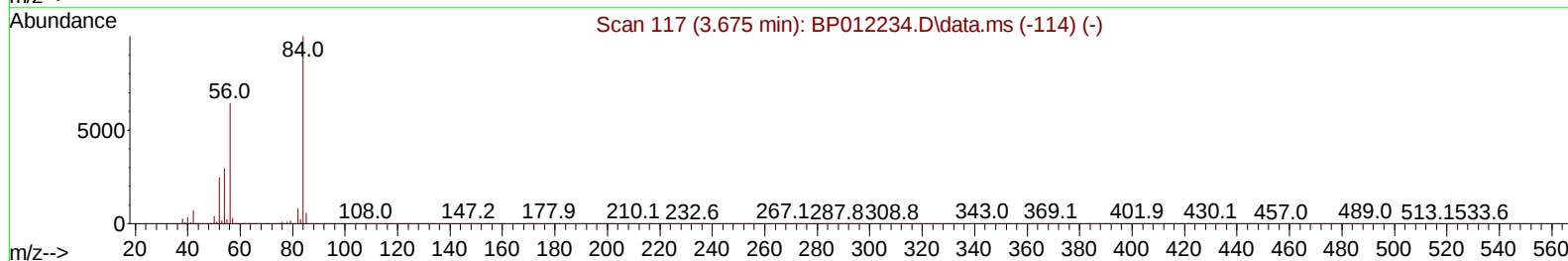
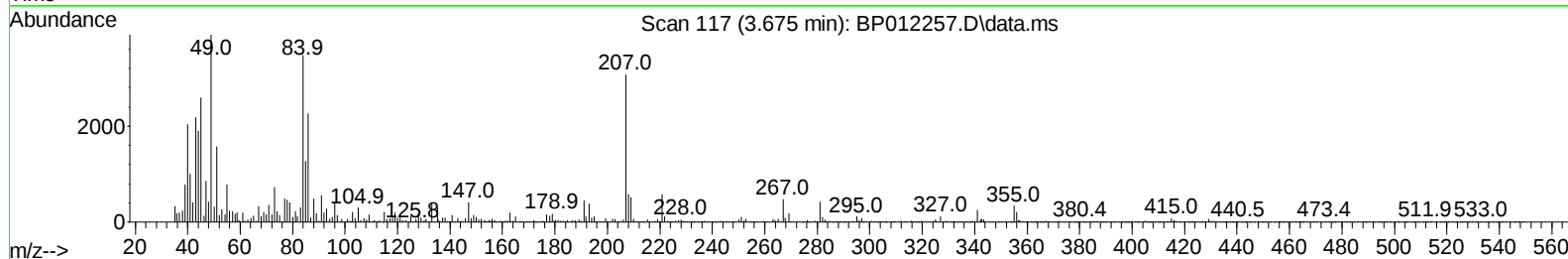
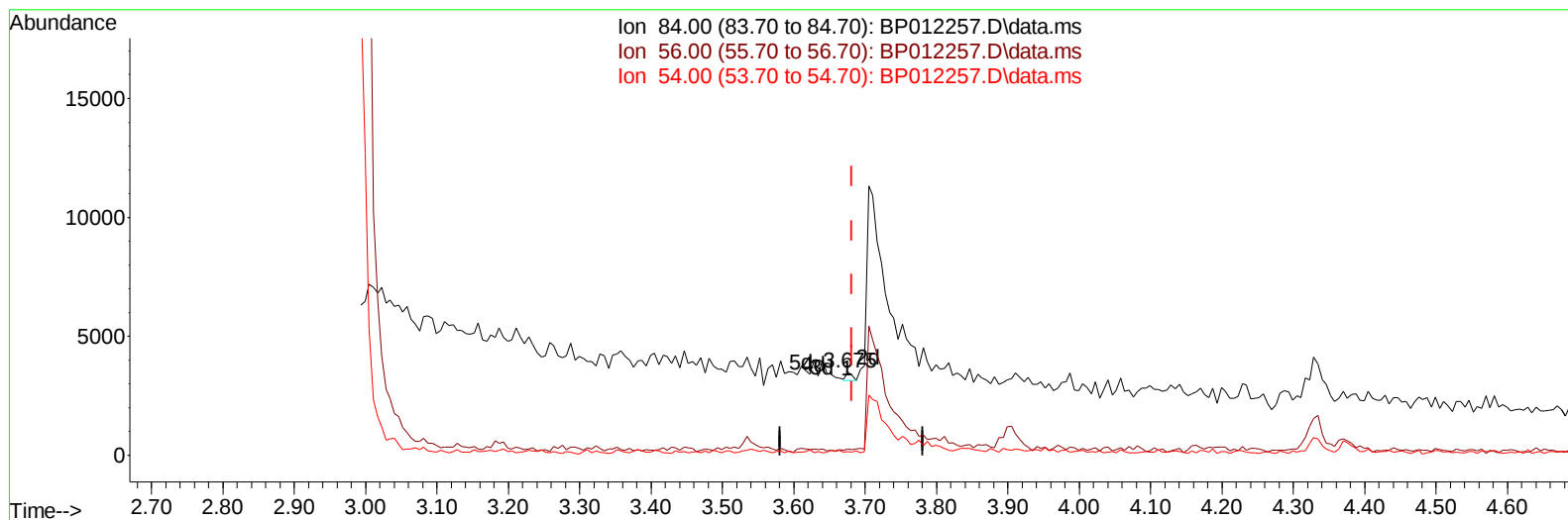
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102022\
 Data File : BP012257.D
 Acq On : 21 Oct 2022 21:49
 Operator : CG/JU
 Sample : N5064-10
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BP6

Manual IntegrationsAPPROVED

Quant Time: Oct 21 22:35:17 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: BP012257.D\data.ms

(4) Pyridine-d5 (S)

3.675min (-0.006) 0.01 ng/u1

response	203	
Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	6.99#
54.00	29.70	4.43#
0.00	0.00	0.00

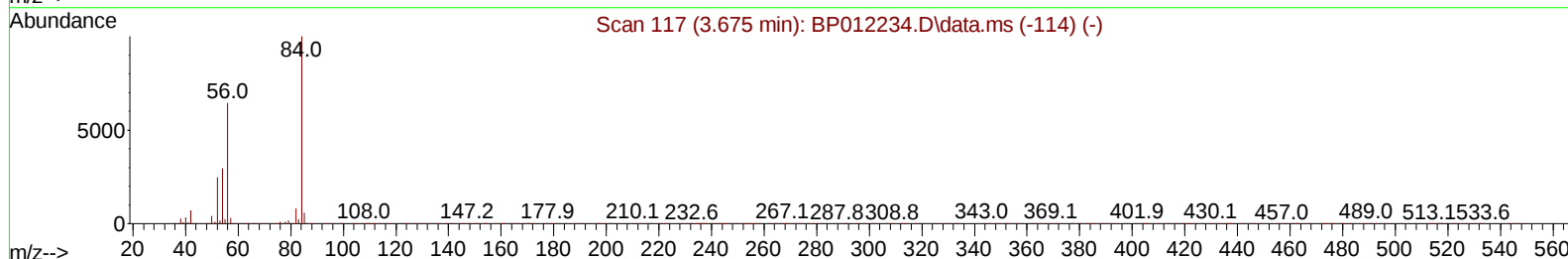
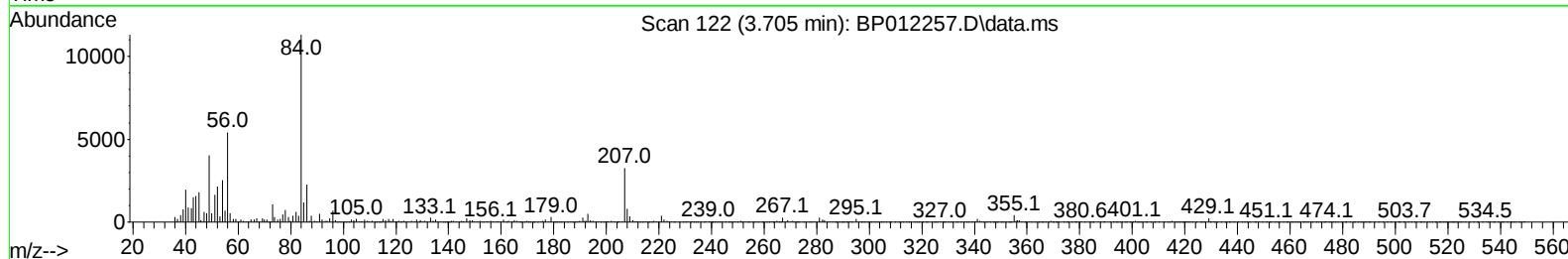
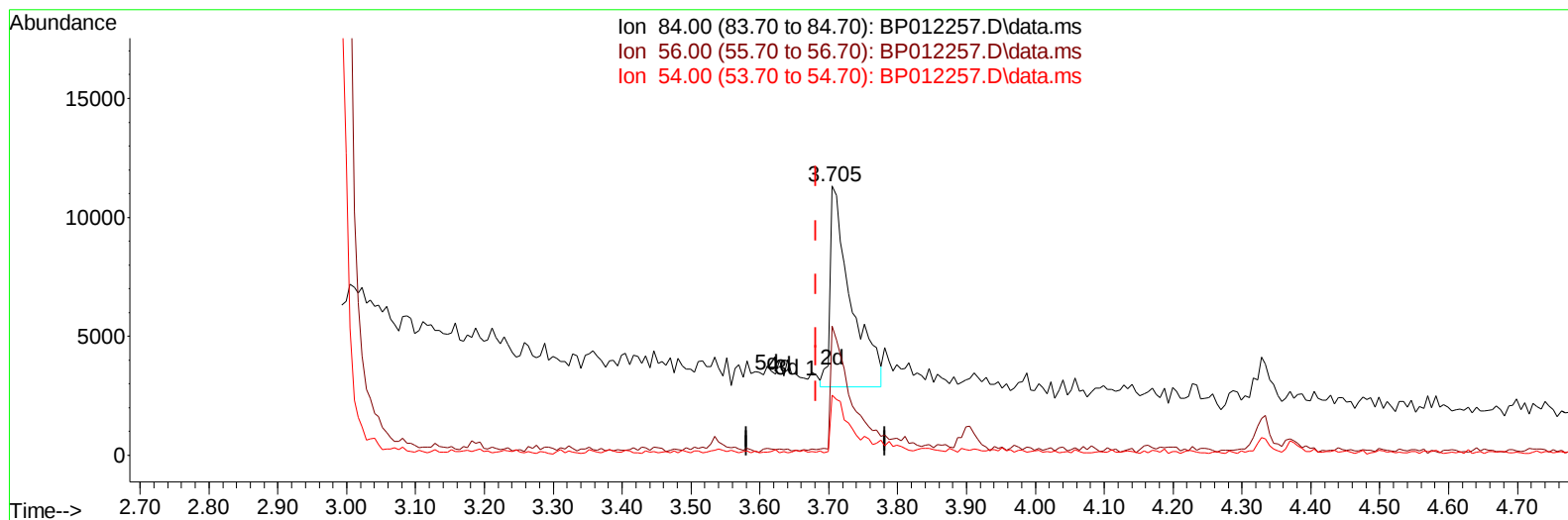
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(4) Pyridine-d5 (S)

3.705min (+ 0.024) 0.66 ng/ul m

response 17817

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	63.60	47.91#
54.00	29.70	22.37#
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	376935	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	1488933	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.469	164	820232	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.234	188	1543977	20.000	ng/ul	0.00
79) Chrysene-d12	21.322	240	1615541	20.000	ng/ul	-0.01
88) Perylene-d12	23.727	264	1753346	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	43545	4.675	ng/uL	0.00
4) Pyridine-d5	3.705	84	17817m	0.664	ng/ul	0.02
7) Phenol-d5	6.993	99	758755	23.317	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	487799	24.717	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	624227	25.201	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	510088	20.169	ng/ul	-0.01
21) Nitrobenzene-d5	8.993	128	307265	27.879	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	314693	26.465	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.246	165	524395	23.511	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.769	131	639787	20.189	ng/ul	-0.01
46) Dimethylphthalate-d6	13.887	166	1484327	26.015	ng/ul	0.00
49) Acenaphthylene-d8	14.163	160	1750913	26.884	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.693	143	164472	15.387	ng/ul	0.00
60) Fluorene-d10	15.469	176	1272778	25.930	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	103471	11.780	ng/ul	0.00
73) Anthracene-d10	17.328	188	1878668	27.759	ng/ul	-0.01
81) Pyrene-d10	19.569	212	2253527	24.500	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.569	264	2423877	27.567	ng/ul	-0.02
Target Compounds						
					Qvalue	
8) Phenol	7.016	94	46416	1.363	ng/ul#	91

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

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