

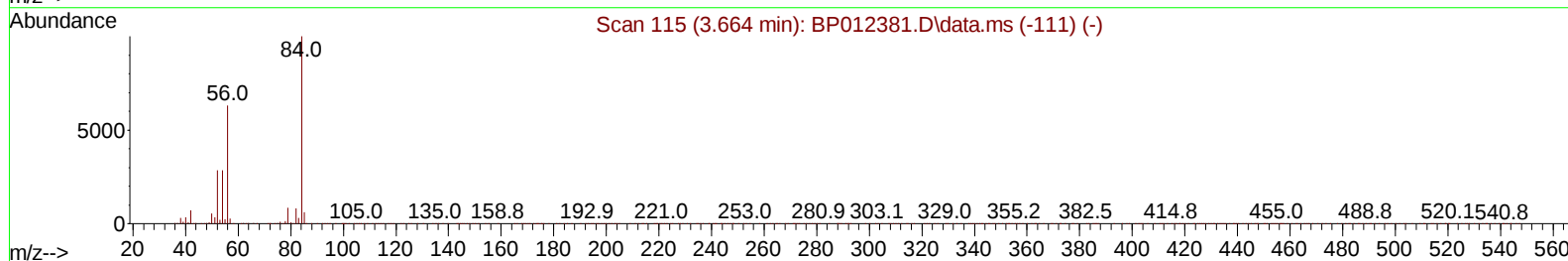
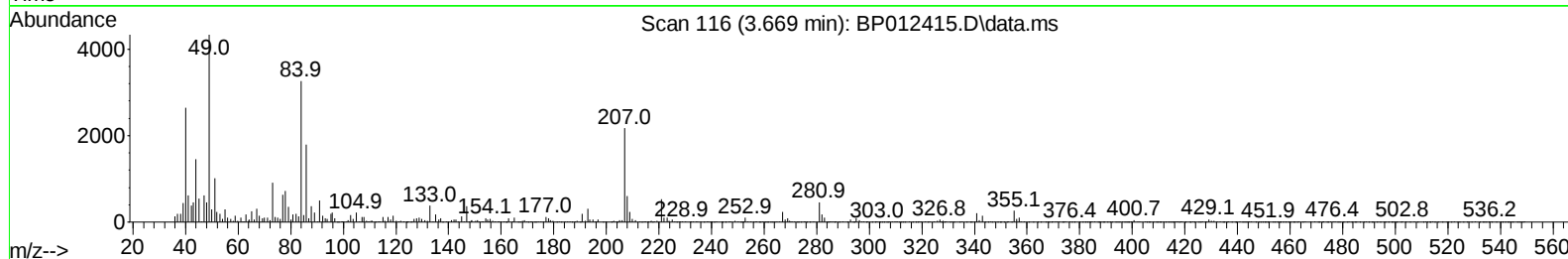
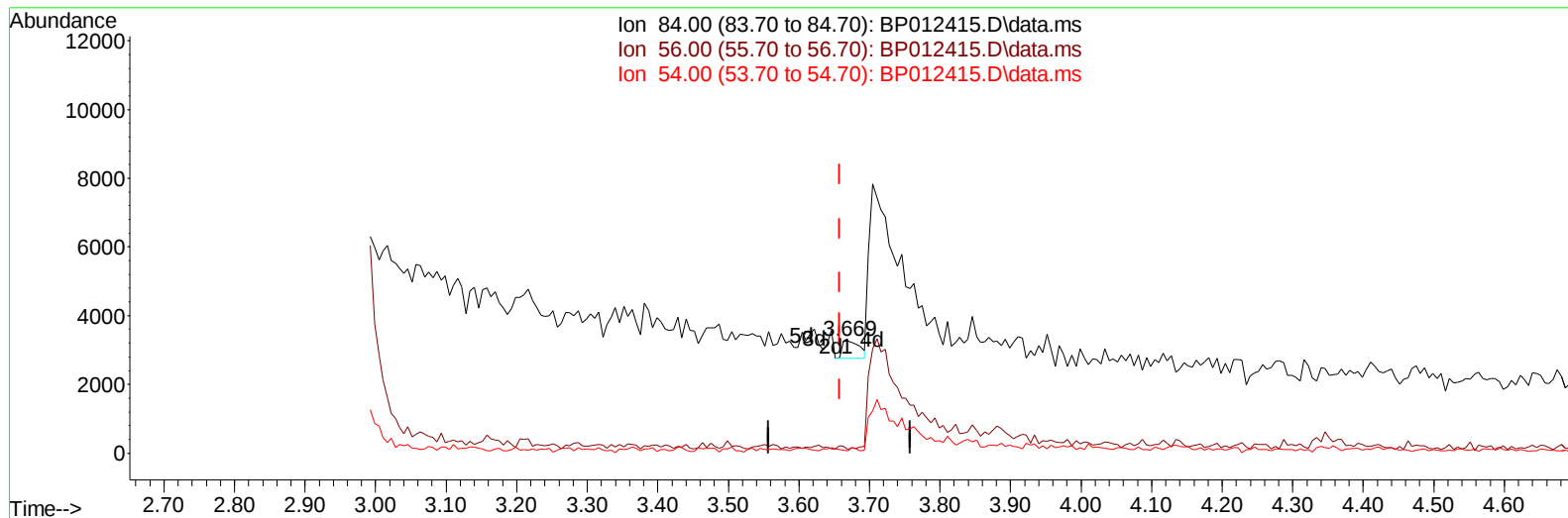
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103122\
Data File : BP012415.D
Acq On : 01 Nov 2022 08:07
Operator : CG/JU
Sample : N5216-10
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BHA87

Manual IntegrationsAPPROVED

Quant Time: Nov 01 08:36:11 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 01 01:03:58 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/01/2022
Supervised By :mohammad ahmed 11/02/2022



TIC: BP012415.D\data.ms

(4) Pyridine-d5 (S)

3.669min (+ 0.011) 0.03 ng/u1

response 838

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	2.95#
54.00	29.80	2.15#
0.00	0.00	0.00

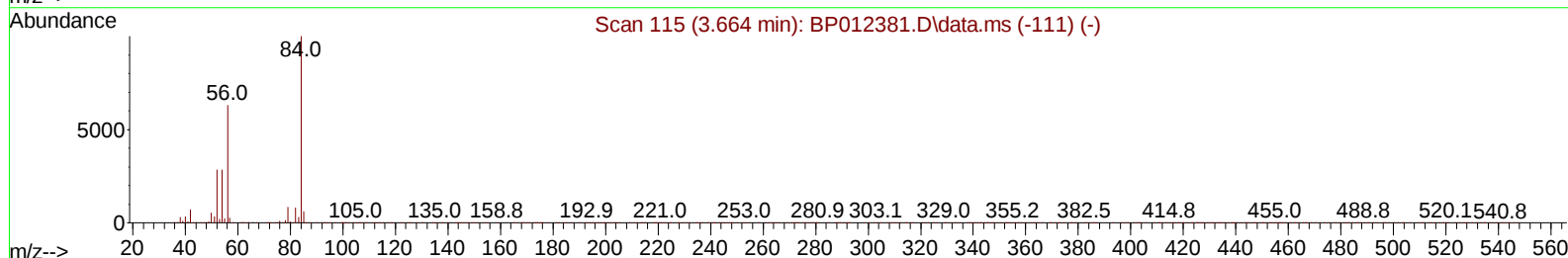
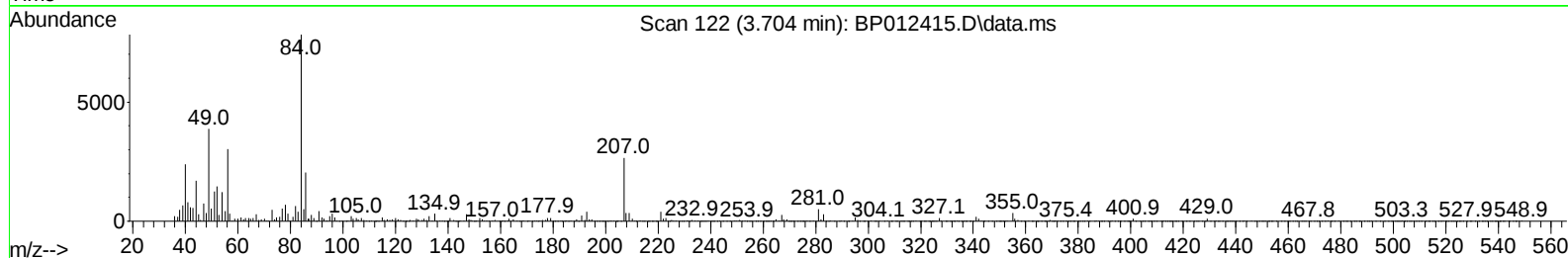
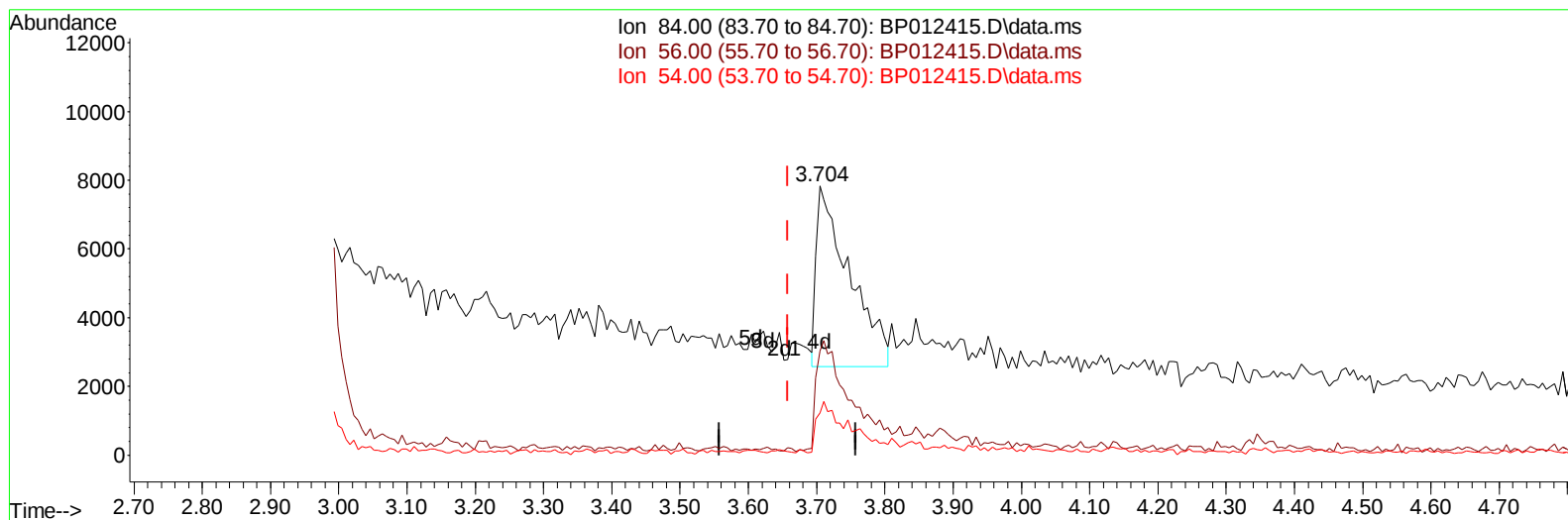
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TIC: BP012415.D\data.ms

(4) Pyridine-d5 (S)

3.704min (+ 0.047) 0.70 ng/ul m

response 17766

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.20	38.78#
54.00	29.80	15.71#
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	7.793	152	357265	20.000	ng/ul	0.00
20) Naphthalene-d8	10.592	136	1610006	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	1052939	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	2324316	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	2440895	20.000	ng/ul	0.00
88) Perylene-d12	23.692	264	2387409	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	31230	3.520	ng/uL	0.00
4) Pyridine-d5	3.704	84	17766m	0.697	ng/ul	0.05
7) Phenol-d5	6.963	99	147933	4.617	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	454592	23.764	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	450723	19.261	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	293311	11.573	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	302526	28.976	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	323558	29.858	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	553611	23.573	ng/ul	0.00
31) 4-Chloroaniline-d4	10.745	131	154984	4.432	ng/ul	0.00
46) Dimethylphthalate-d6	13.857	166	2118172	29.387	ng/ul	0.00
49) Acenaphthylene-d8	14.133	160	2269618	27.554	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	74131	5.434	ng/ul	0.00
60) Fluorene-d10	15.439	176	1913788	31.011	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.574	200	394084	32.565	ng/ul	0.00
73) Anthracene-d10	17.304	188	3505776	35.005	ng/ul	0.00
81) Pyrene-d10	19.545	212	4286351	34.977	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.539	264	4145460	35.537	ng/ul	0.00

Target Compounds Qvalue

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

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