

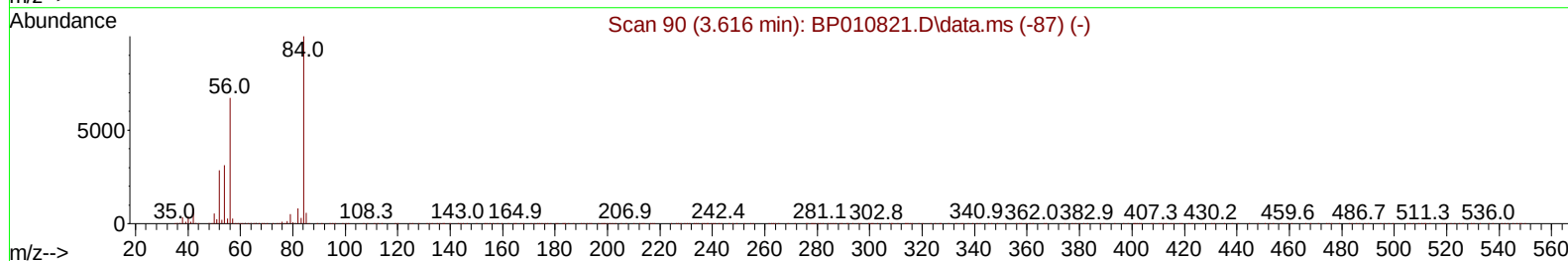
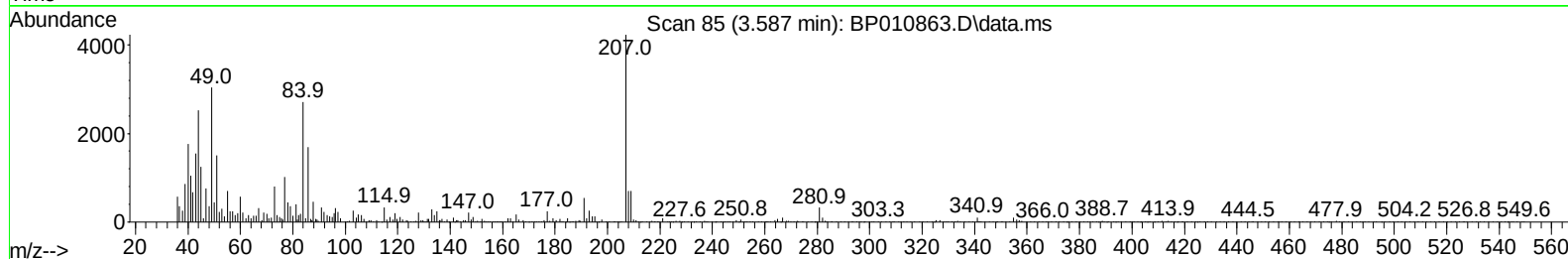
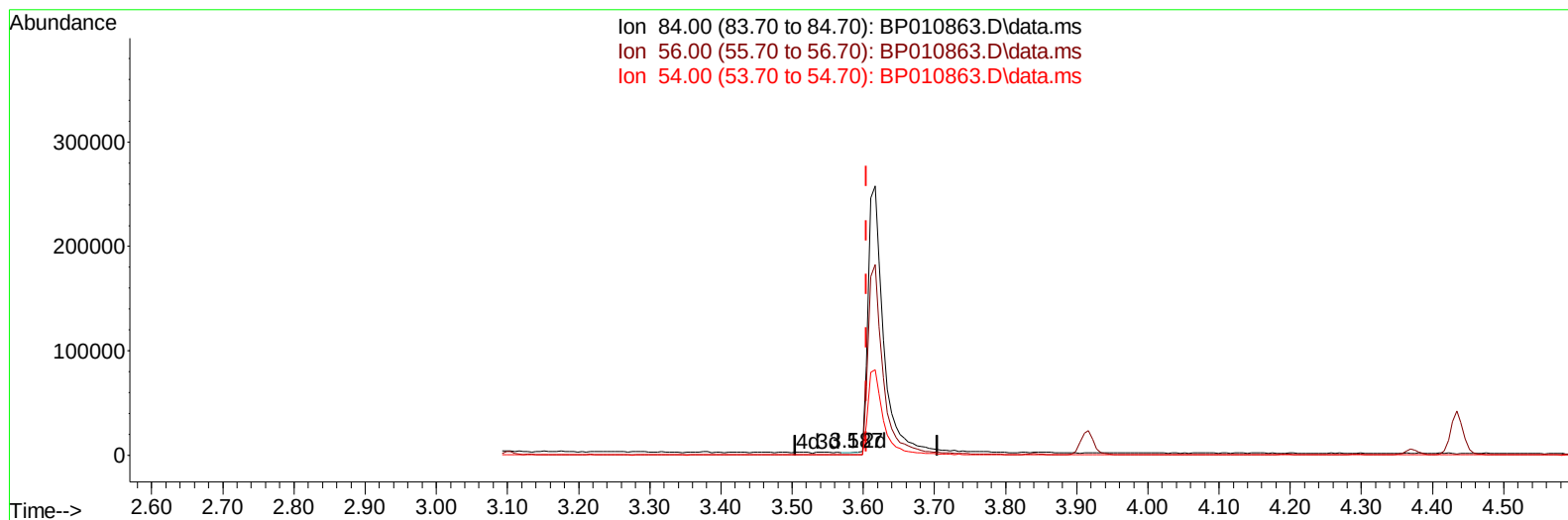
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010863.D
 Acq On : 26 Jun 2022 11:53
 Operator : CG/JU
 Sample : N3392-04 10X
 Misc :
 ALS Vial : 45 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEX9

Manual IntegrationsAPPROVED

Quant Time: Jun 26 22:58:38 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 24 22:42:26 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/27/2022
 Supervised By :mohammad ahmed 07/01/2022



TIC: BP010863.D\data.ms

(4) Pyridine-d5 (S)

3.587min (-0.018) 0.01 ng/ul

response 392

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	8.88#
54.00	44.80	4.86#
0.00	0.00	0.00

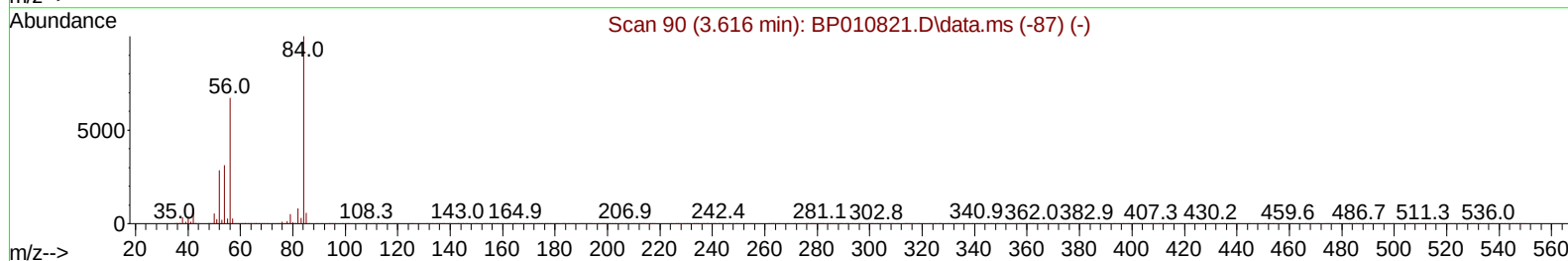
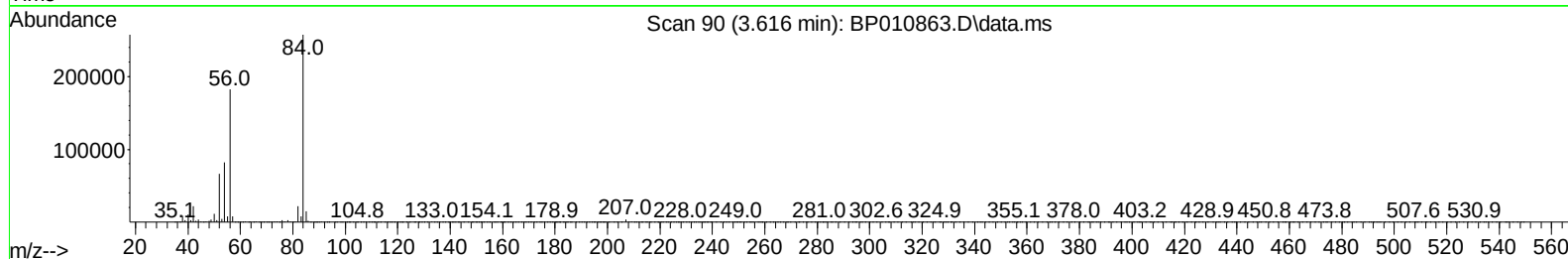
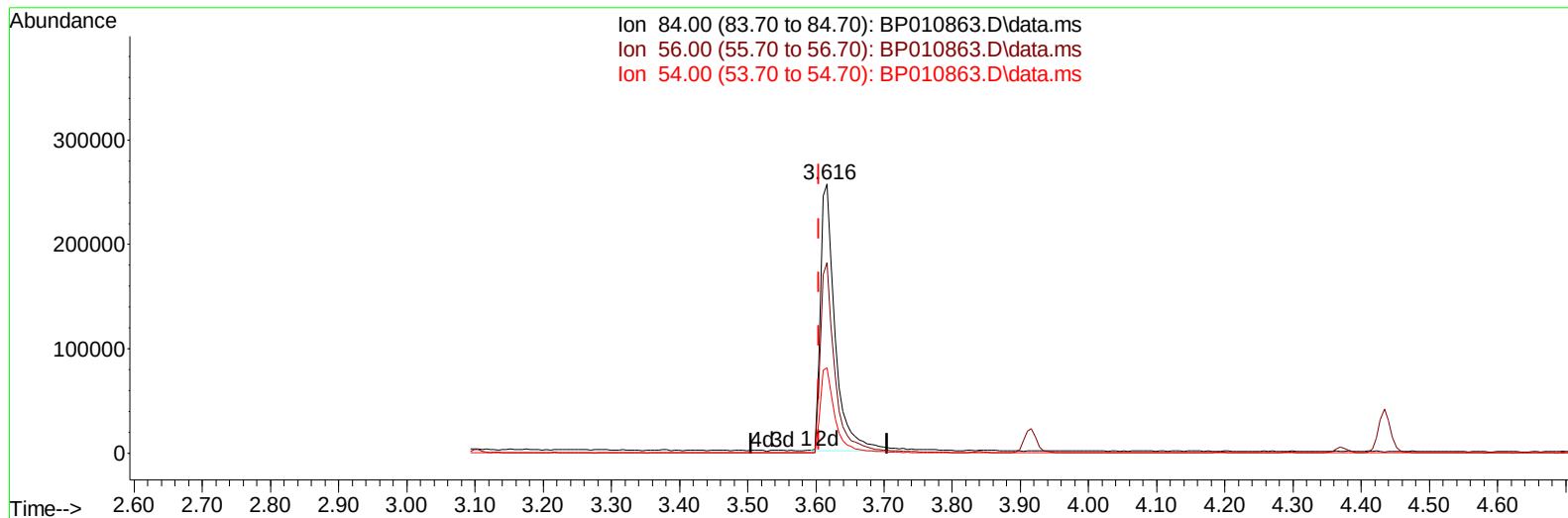
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TIC: BP010863.D\data.ms

(4) Pyridine-d5 (S)

3.616min (+ 0.012) 10.99 ng/ul m

response 383207

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	70.78
54.00	44.80	31.71#
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	7.710	152	495550	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.504	136	1996100	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.369	164	1052607	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.133	188	1885101	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.245	240	1744966	20.000	ng/ul	#-0.01
88) Perylene-d12	23.609	264	1852284	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	30571	2.334	ng/uL	0.00
4) Pyridine-d5	3.616	84	383207m	10.989	ng/ul	0.01
7) Phenol-d5	6.887	99	245633	6.262	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.051	67	875435	34.124	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.246	132	839452	25.793	ng/ul	0.00
15) 4-Methylphenol-d8	8.428	113	486314	15.682	ng/ul	0.00
21) Nitrobenzene-d5	8.875	128	525936	41.389	ng/ul	0.00
24) 2-Nitrophenol-d4	9.592	143	472946	44.987	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.128	165	800482	30.816	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.651	131	1192076	28.114	ng/ul	0.00
46) Dimethylphthalate-d6	13.786	166	2736352	37.950	ng/ul	-0.01
49) Acenaphthylene-d8	14.057	160	3259268	35.417	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.580	143	54805	4.866	ng/ul	0.00
60) Fluorene-d10	15.369	176	2321369	37.516	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	182812	29.271	ng/ul	-0.01
73) Anthracene-d10	17.233	188	3437290	40.983	ng/ul	-0.01
81) Pyrene-d10	19.480	212	3689360	38.808	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.456	264	3879423	42.383	ng/ul	-0.02

Target Compounds Qvalue

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

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