

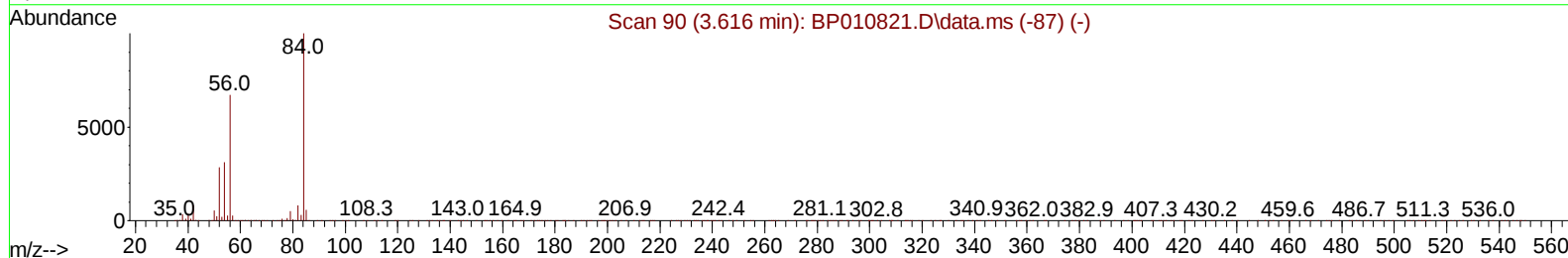
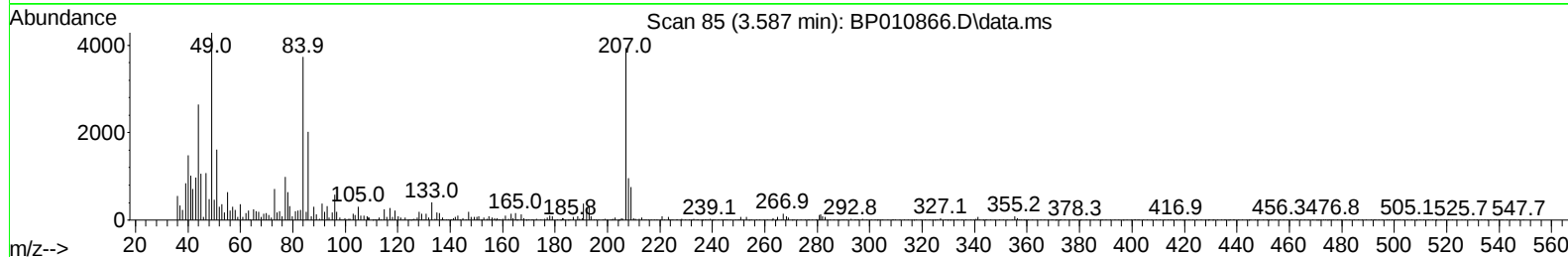
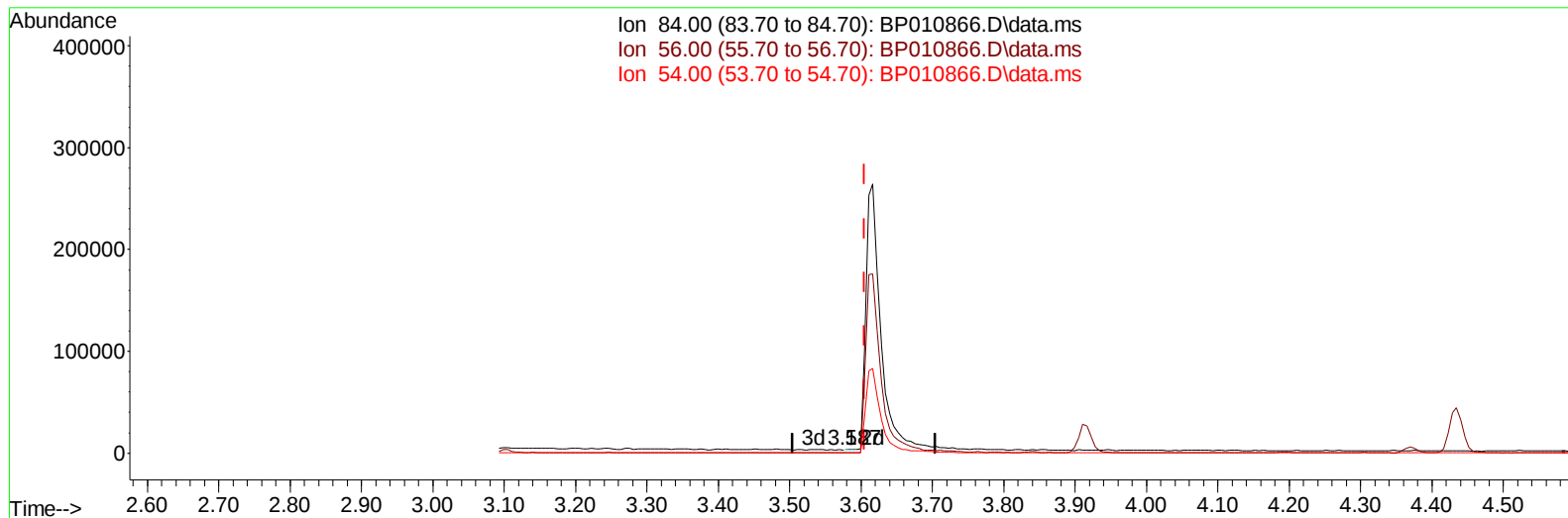
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010866.D
 Acq On : 26 Jun 2022 13:34
 Operator : CG/JU
 Sample : N3392-05 10X
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY0

Manual IntegrationsAPPROVED

Quant Time: Jun 26 22:59:05 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: BP010866.D\data.ms

(4) Pyridine-d5 (S)

3.587min (-0.018) 0.02 ng/u1

response 810

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	5.96#
54.00	44.80	4.65#
0.00	0.00	0.00

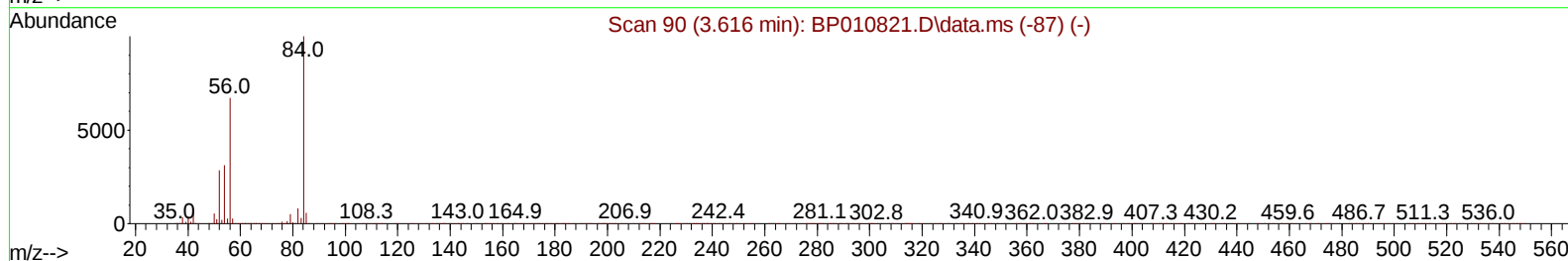
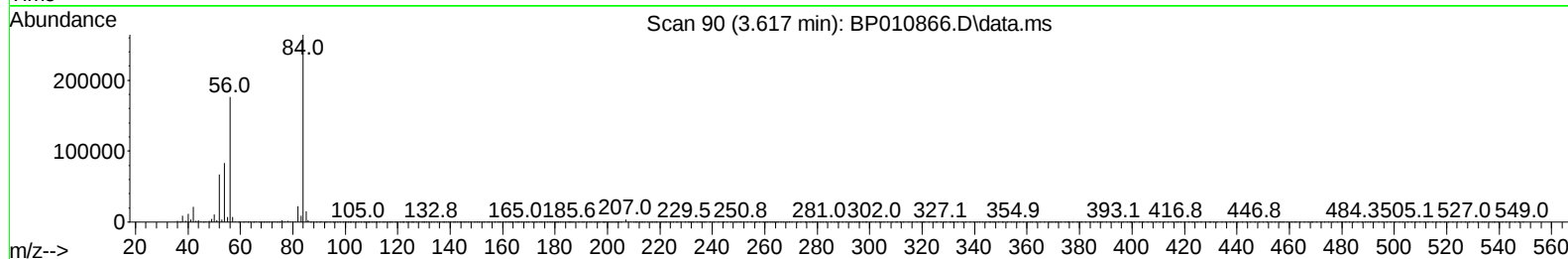
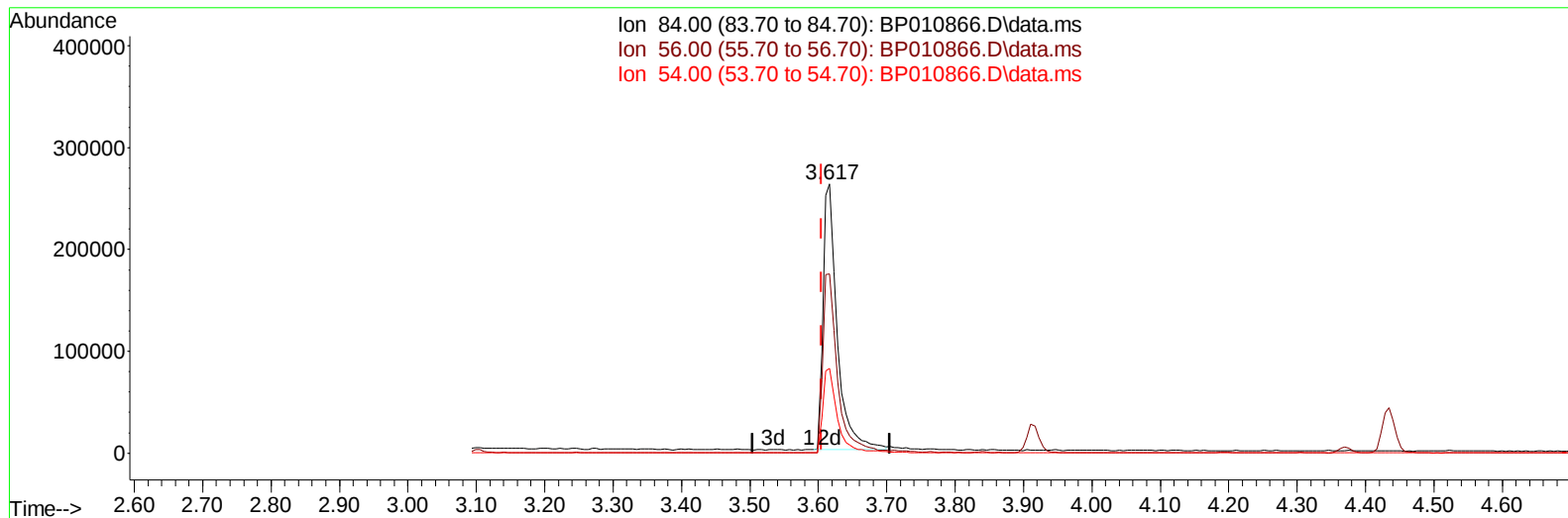
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(4) Pyridine-d5 (S)

3.617min (+ 0.012) 10.71 ng/ul m

response 376728

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	66.71#
54.00	44.80	31.41#
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	7.711	152	499976	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.505	136	2034235	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.369	164	1072974	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.134	188	1919469	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.245	240	1764625	20.000	ng/ul	#-0.01
88) Perylene-d12	23.610	264	1908567	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	32200	2.436	ng/uL	0.00
4) Pyridine-d5	3.617	84	376728m	10.708	ng/ul	0.01
7) Phenol-d5	6.887	99	283407	7.161	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	893662	34.526	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.246	132	886445	26.995	ng/ul	0.00
15) 4-Methylphenol-d8	8.428	113	540973	17.291	ng/ul	0.00
21) Nitrobenzene-d5	8.875	128	539580	41.667	ng/ul	0.00
24) 2-Nitrophenol-d4	9.593	143	490792	45.810	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.128	165	825107	31.169	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.652	131	1200119	27.773	ng/ul	0.00
46) Dimethylphthalate-d6	13.787	166	2706453	36.823	ng/ul	-0.01
49) Acenaphthylene-d8	14.057	160	3252667	34.675	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.581	143	63720	5.550	ng/ul	0.00
60) Fluorene-d10	15.369	176	2291945	36.337	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.487	200	183224	28.812	ng/ul	-0.01
73) Anthracene-d10	17.234	188	3376324	39.535	ng/ul	-0.01
81) Pyrene-d10	19.481	212	3605496	37.504	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.457	264	3840712	40.723	ng/ul	-0.02

Target Compounds Qvalue

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

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