

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062522\
 Data File : BP010857.D
 Acq On : 26 Jun 2022 08:30
 Operator : CG/JU
 Sample : N3392-14 10X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXEY9

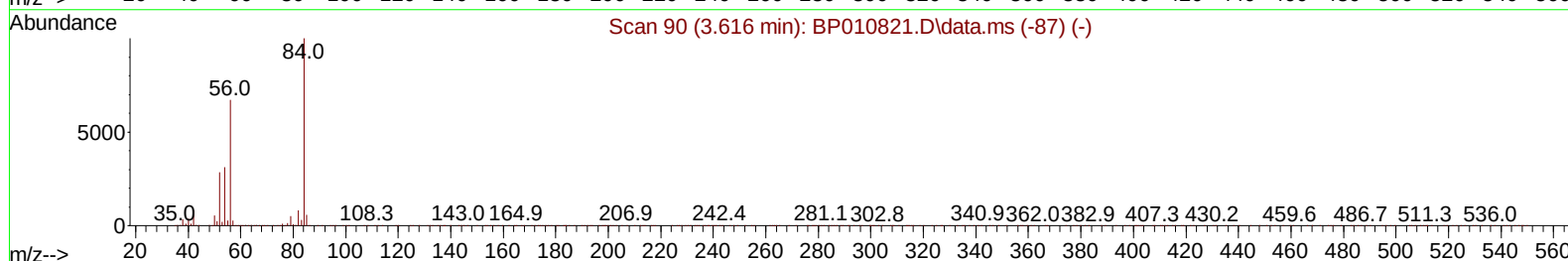
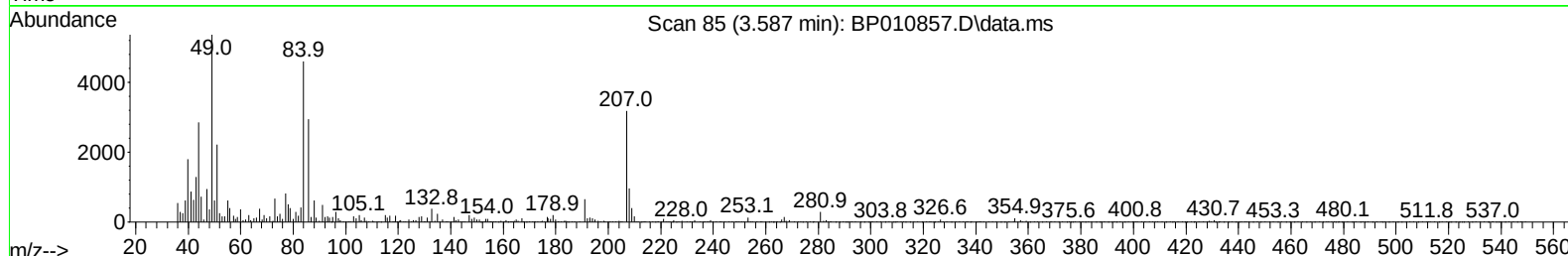
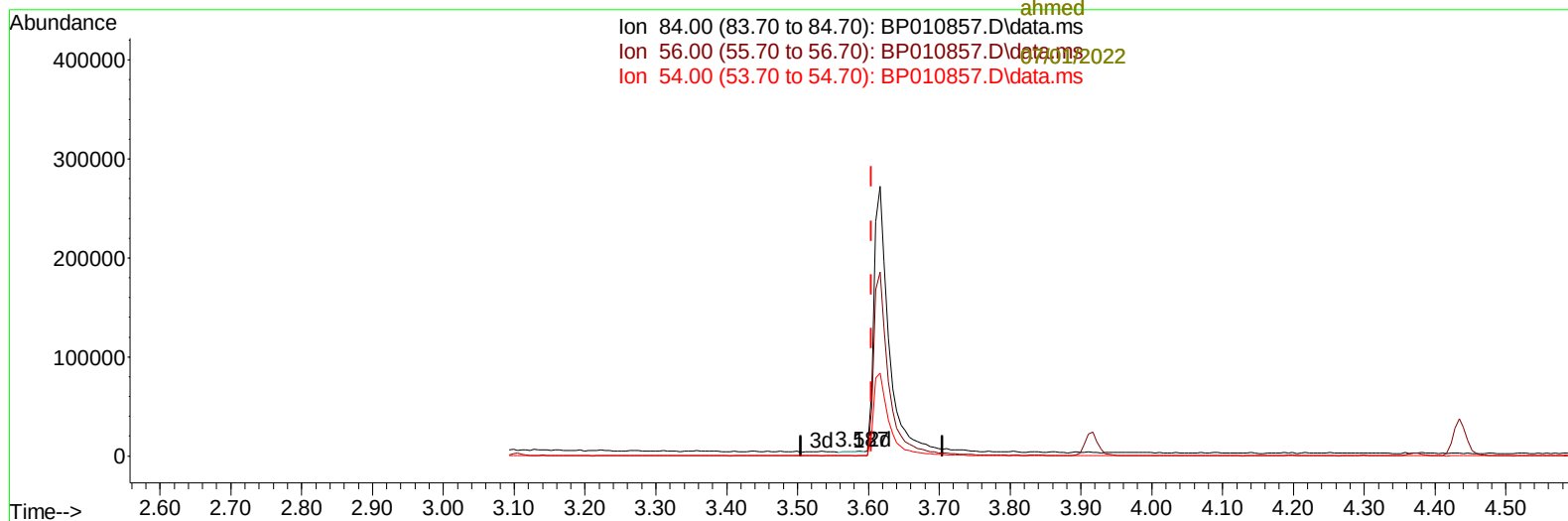
Manual IntegrationsAPPROVED

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

06/27/2022
 Supervised By :mohammad
 ahmed

Quant Time: Jun 26 22:57:47 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

Ion 84.00 (83.70 to 84.70): BP010857.D\data.ms
 Ion 56.00 (55.70 to 56.70): BP010857.D\data.ms
 Ion 54.00 (53.70 to 54.70): BP010857.D\data.ms



TIC: BP010857.D\data.ms

(4) Pyridine-d5 (S)

3.587min (-0.018) 0.03 ng/ul

response 1028

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	8.86#
54.00	44.80	3.63#
0.00	0.00	0.00

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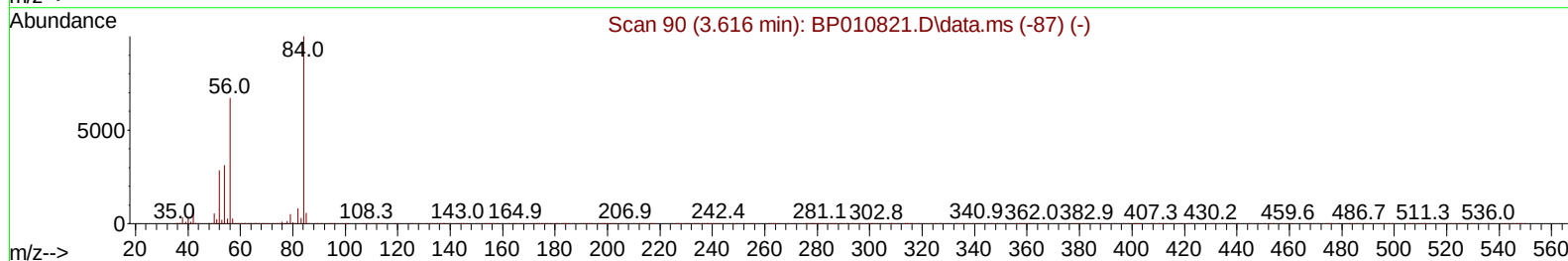
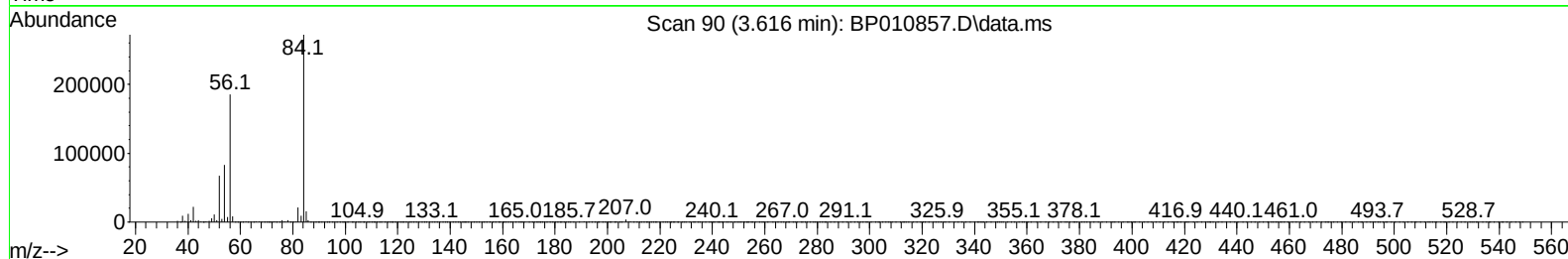
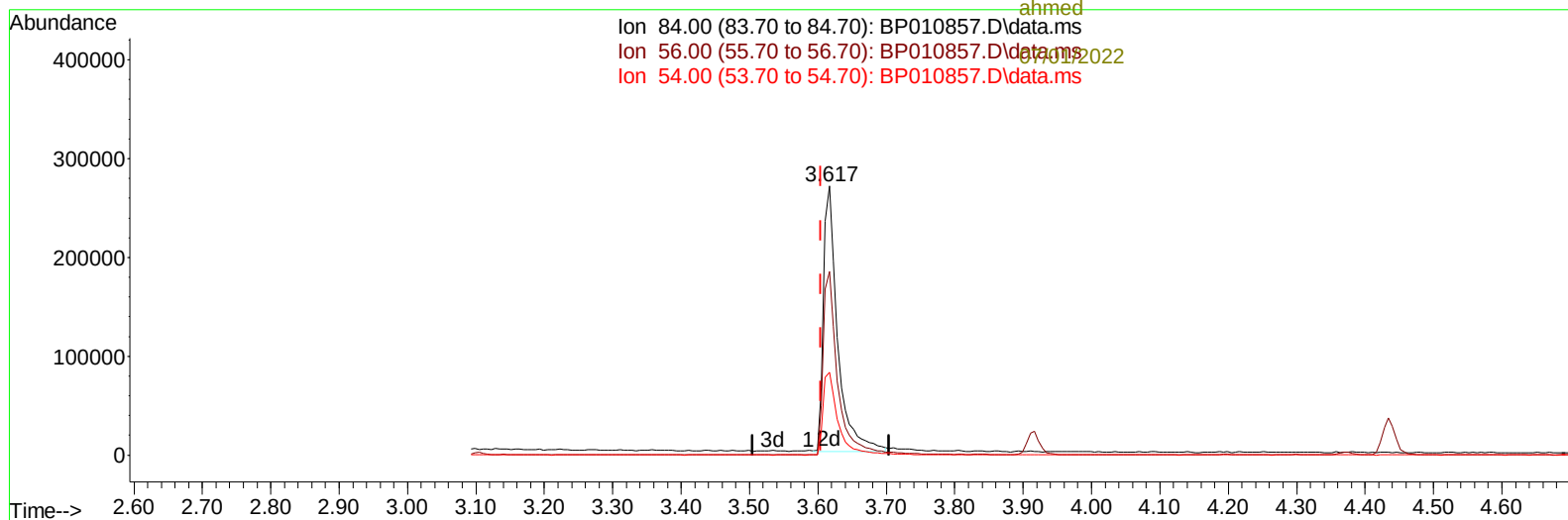
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 Ion 56.00 (55.70 to 56.70): BP010857.D\data.ms
 Ion 54.00 (53.70 to 54.70): BP010857.D\data.ms



TIC: BP010857.D\data.ms

(4) Pyridine-d5 (S)

3.616min (+ 0.012) 11.43 ng/ul m

response 390059

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	84.50	68.16
54.00	44.80	30.68#
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----07/01/2022						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.710	152	484898	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.510	136	1949625	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.369	164	1044550	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.133	188	2043390	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.245	240	1994262	20.000	ng/ul	#-0.01
88) Perylene-d12	23.610	264	2091417	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	33136	2.585	ng/uL	0.00
4) Pyridine-d5	3.616	84	390059m	11.431	ng/ul	0.01
7) Phenol-d5	6.893	99	286620	7.467	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.052	67	843320	33.594	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.246	132	862407	27.080	ng/ul	0.00
15) 4-Methylphenol-d8	8.428	113	526683	17.357	ng/ul	0.00
21) Nitrobenzene-d5	8.875	128	503825	40.594	ng/ul	0.00
24) 2-Nitrophenol-d4	9.593	143	458513	44.654	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.134	165	789442	31.116	ng/ul	0.00
31) 4-Chloroaniline-d4	10.651	131	1207947	29.167	ng/ul	0.00
46) Dimethylphthalate-d6	13.787	166	2686807	37.550	ng/ul	-0.01
49) Acenaphthylene-d8	14.057	160	3134896	34.328	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.581	143	83931	7.510	ng/ul	0.00
60) Fluorene-d10	15.369	176	2342050	38.142	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.486	200	212374	31.370	ng/ul	-0.01
73) Anthracene-d10	17.233	188	3812071	41.931	ng/ul	-0.01
81) Pyrene-d10	19.480	212	4274439	39.342	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.457	264	4459925	43.154	ng/ul	-0.02

Target Compounds Qvalue

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

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