

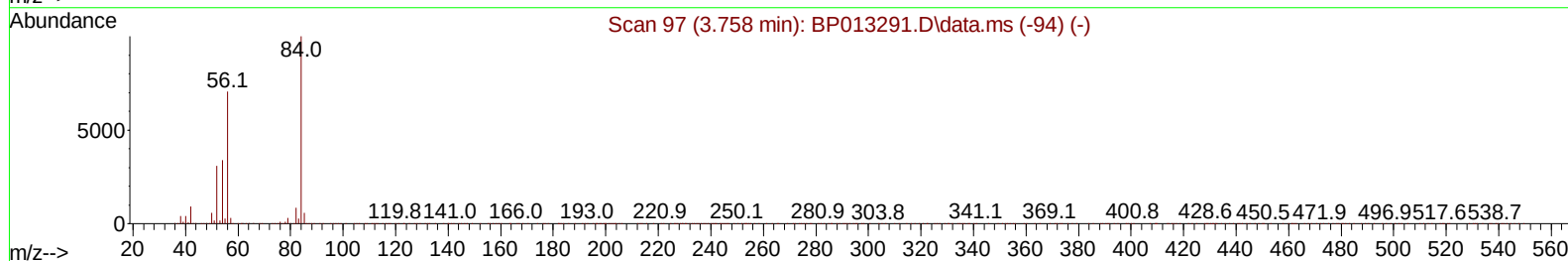
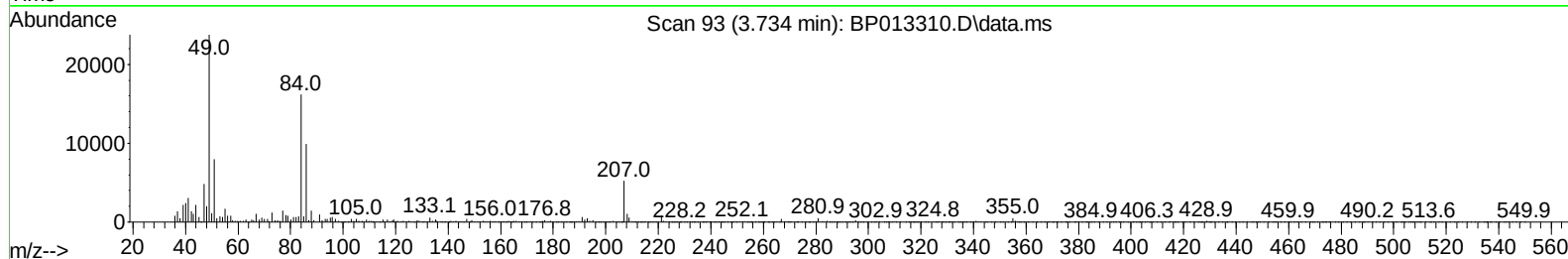
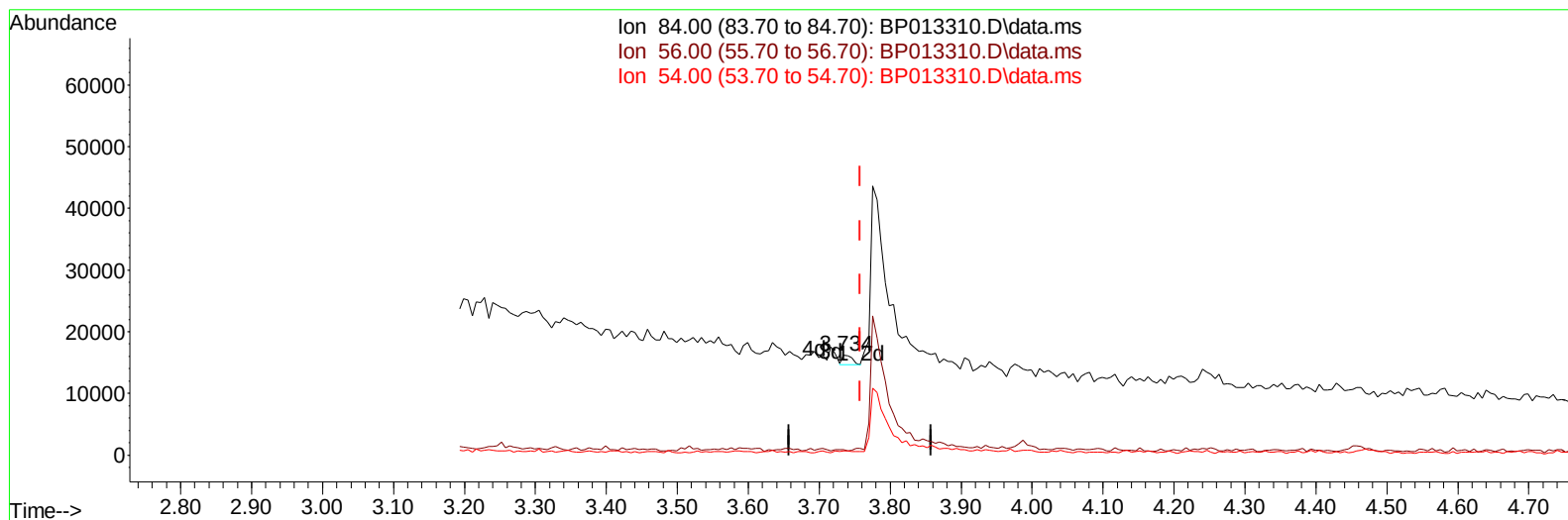
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013310.D
Acq On : 24 Dec 2022 21:28
Operator : CG/JU
Sample : N6016-14
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD6

Manual IntegrationsAPPROVED

Quant Time: Dec 25 00:25:03 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Dec 24 00:19:17 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 12/25/2022
Supervised By :Yogesh Patel 12/25/2022



TIC: BP013310.D\data.ms

(4) Pyridine-d5 (S)

3.734min (-0.023) 0.04 ng/uI

response 1573

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	5.18#
54.00	30.50	4.08#
0.00	0.00	0.00

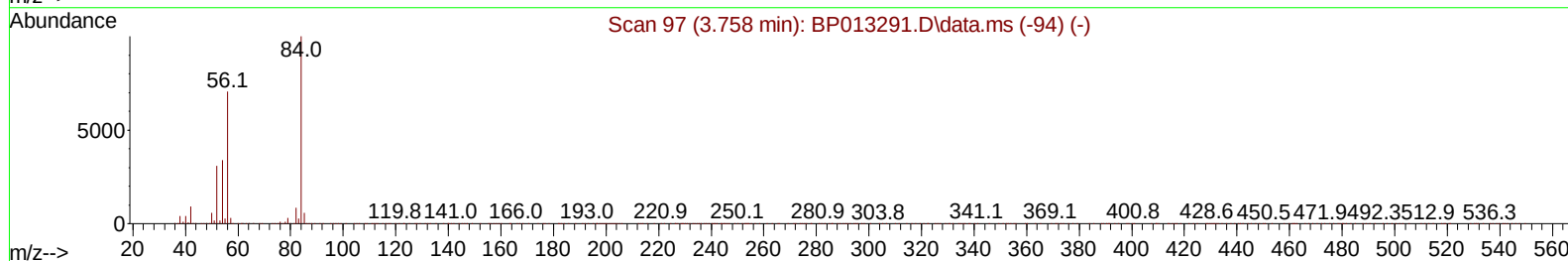
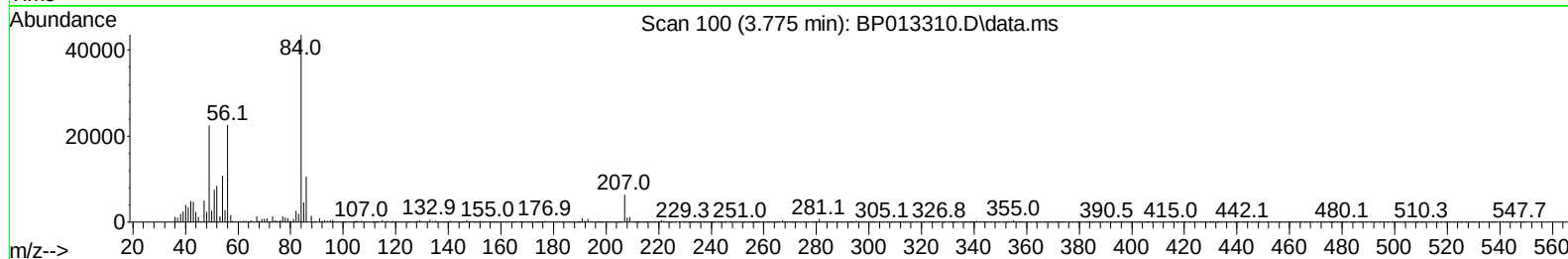
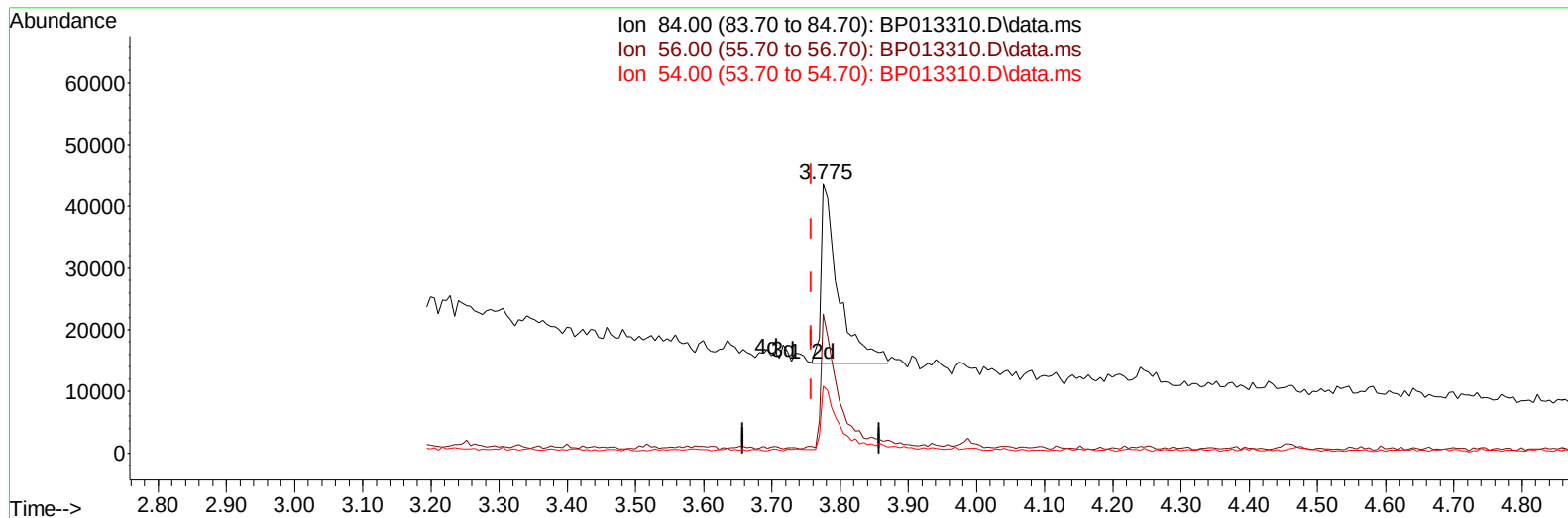
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TIC: BP013310.D\data.ms

(4) Pyridine-d5 (S)

3.775min (+ 0.018) 1.19 ng/ul m

response 52123

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.30	51.77#
54.00	30.50	24.90
0.00	0.00	0.00

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 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	659855	20.000	ng/ul	0.00
20) Naphthalene-d8	10.752	136	2569859	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.581	164	1321082	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.345	188	2501477	20.000	ng/ul	0.00
79) Chrysene-d12	21.433	240	2667945	20.000	ng/ul	0.00
88) Perylene-d12	23.916	264	3090905	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.328	96	47096	3.052	ng/uL	0.00
4) Pyridine-d5	3.775	84	52123m	1.189	ng/ul	0.02
7) Phenol-d5	7.099	99	826018	15.369	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.263	67	587957	18.134	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	720038	17.018	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113	419770	9.532	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128	348808	18.474	ng/ul	0.00
24) 2-Nitrophenol-d4	9.834	143	385071	18.115	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165	660299	15.992	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.893	131	609624	10.459	ng/ul	0.00
46) Dimethylphthalate-d6	13.987	166	1893373	18.648	ng/ul	-0.01
49) Acenaphthylene-d8	14.275	160	2067412	18.531	ng/ul	0.00
54) 4-Nitrophenol-d4	14.787	143	160348	8.780	ng/ul	0.00
60) Fluorene-d10	15.575	176	1496251	17.419	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.698	200	128356	7.974	ng/ul	0.00
73) Anthracene-d10	17.439	188	2062827	18.279	ng/ul	0.00
81) Pyrene-d10	19.675	212	2843246	18.566	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.751	264	2957117	19.194	ng/ul	0.00

Target Compounds Qvalue

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

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