Quantitation Report (Qedit)

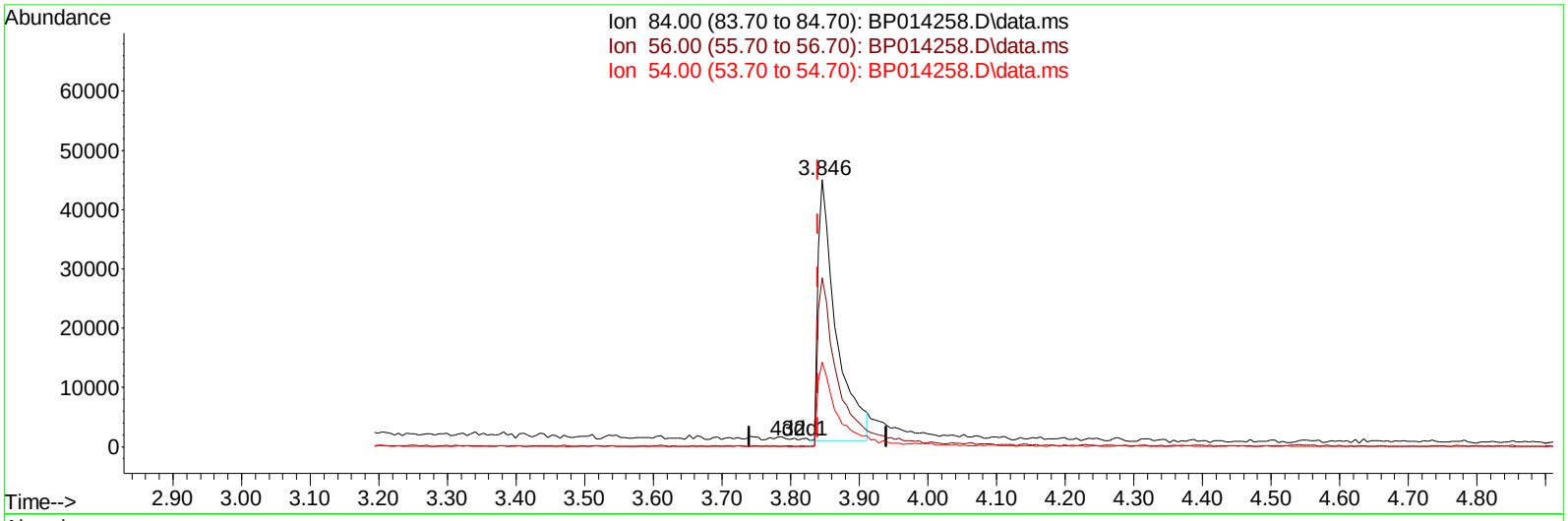
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014258.D
Acq On : 23 Mar 2023 15:16
Operator : CG/JU
Sample : 02010-03
Misc :
ALS Vial : 38 Sample Multiplier: 1

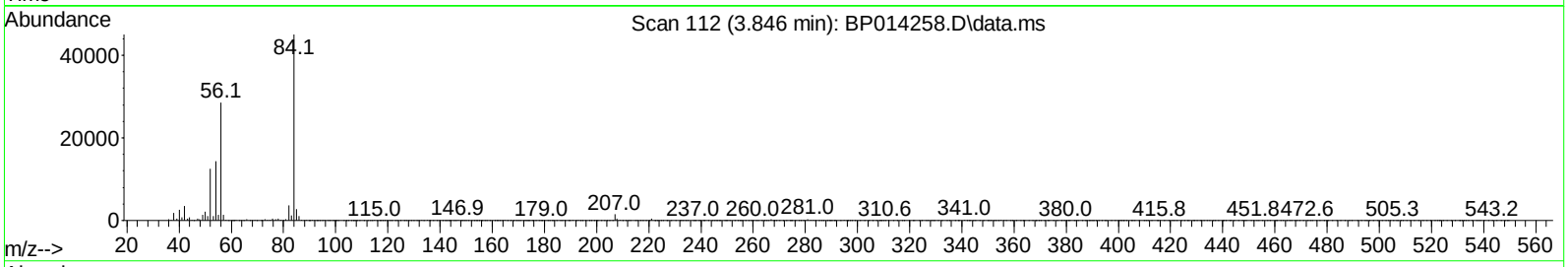
Quant Time: Mar 24 02:08:44 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 23 04:56:06 2023
Response via : Initial Calibration

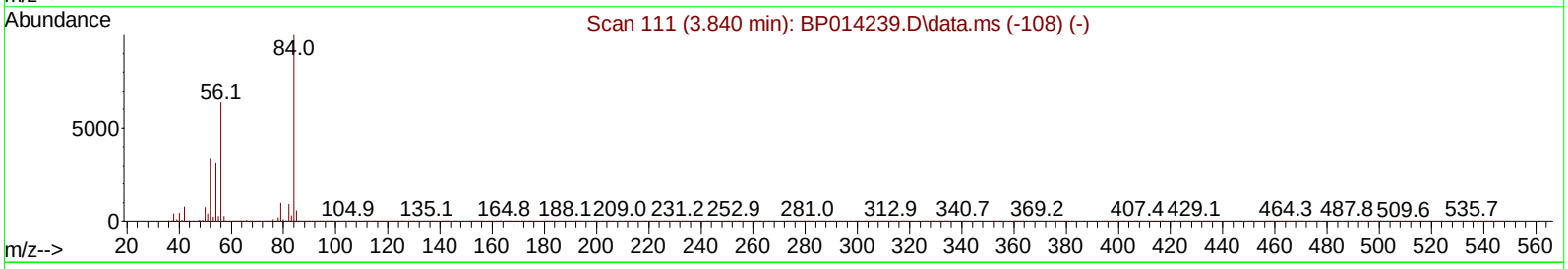
Instrument :
BNA_P
ClientSampleId :
CB949

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/24/2023
Supervised By :Jagrut Upadhyay 03/24/2023







TIC: BP014258.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.006) 7.45 ng/ul

response 80240

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	63.42
54.00	33.20	31.76
0.00	0.00	0.00

Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
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Acq On : 23 Mar 2023 15:16
Operator : CG/JU
Sample : 02010-03
Misc :
ALS Vial : 38 Sample Multiplier: 1

Quant Time: Mar 24 02:08:44 2023
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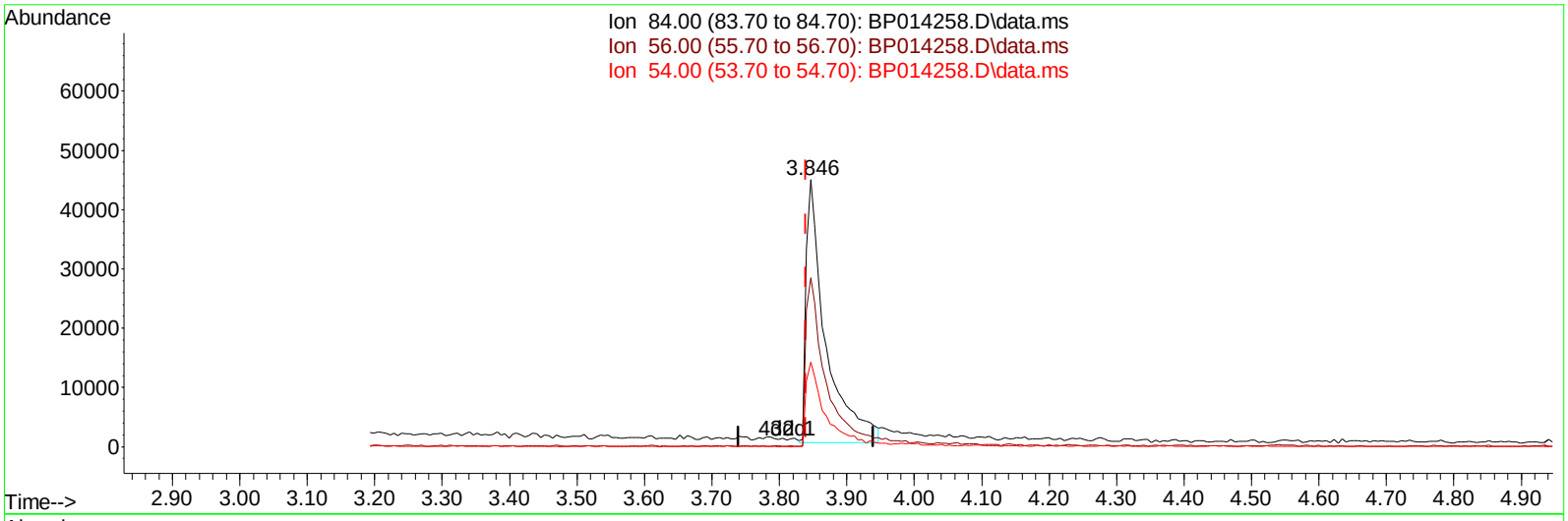

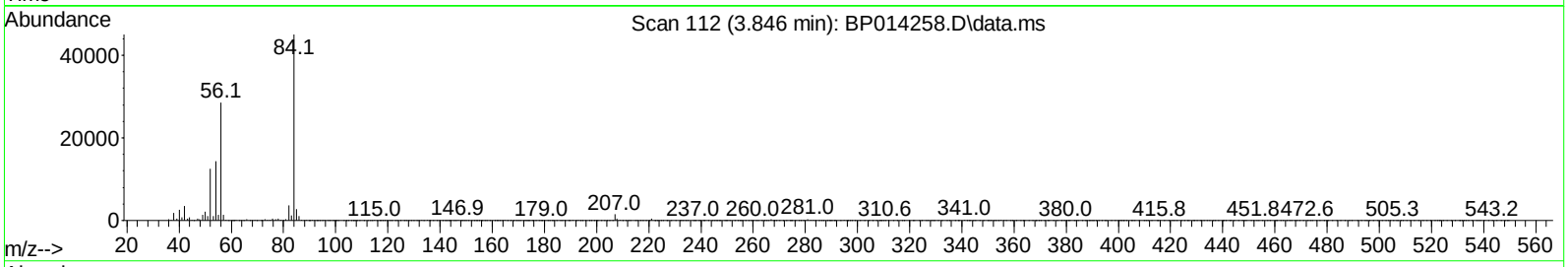

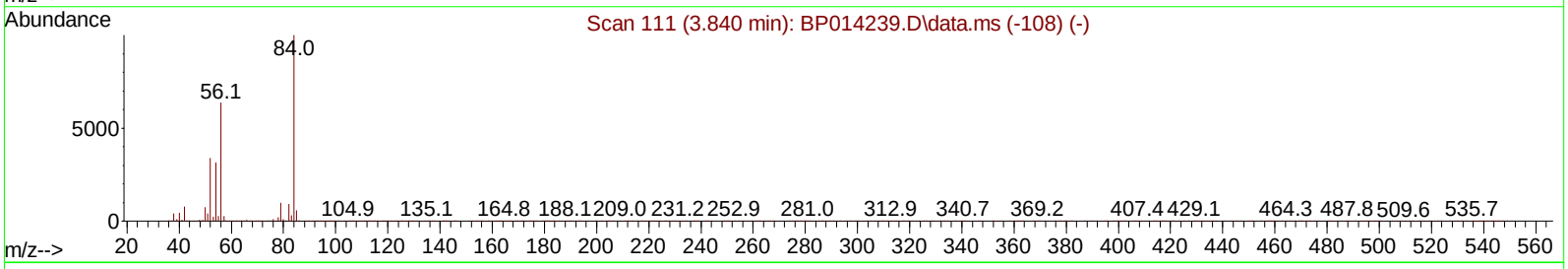

TIC: BP014258.D\data.ms

(4) Pyridine-d5 (S)

3.846min (+ 0.006) 8.24 ng/ul m

response 88789

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	62.80	63.42
54.00	33.20	31.76
0.00	0.00	0.00

Quantitation Report (QT/LSC Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032223\
Data File : BP014258.D
Acq On : 23 Mar 2023 15:16
Operator : CG/JU
Sample : 02010-03
Misc :
ALS Vial : 38 Sample Multiplier: 1

Quant Time: Mar 24 02:08:44 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 23 04:56:06 2023
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/24/2023
Supervised By :Jagrut Upadhyay 03/24/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.034	152	147073	20.000	ng/ul	0.00
20) Naphthalene-d8	10.858	136	675018	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.681	164	492330	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.440	188	1150504	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.522	240	1140556	20.000	ng/ul	# 0.00
88) Perylene-d12	24.039	264	1179639	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.405	96	14972	3.846	ng/uL	0.00
4) Pyridine-d5	3.846	84	88789m	8.238	ng/ul	0.00
7) Phenol-d5	7.199	99	80689	6.217	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.358	67	235795	31.152	ng/ul	0.00
11) 2-Chlorophenol-d4	7.564	132	215082	21.751	ng/ul	0.00
15) 4-Methylphenol-d8	8.752	113	172423	16.264	ng/ul	0.00
21) Nitrobenzene-d5	9.211	128	149257	27.712	ng/ul	0.00
24) 2-Nitrophenol-d4	9.940	143	162903	25.267	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.481	165	315400	25.999	ng/ul	0.00
31) 4-Chloroaniline-d4	10.999	131	372850	23.225	ng/ul	0.00
46) Dimethylphthalate-d6	14.087	166	1301451	31.148	ng/ul	0.00
49) Acenaphthylene-d8	14.375	160	1340517	30.286	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	27142	3.686	ng/ul	0.02
60) Fluorene-d10	15.675	176	1094412	30.906	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200	111359	12.806	ng/ul	0.00
73) Anthracene-d10	17.540	188	1868958	33.463	ng/ul	0.00
81) Pyrene-d10	19.763	212	2284738	36.320	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.874	264	2178077	34.712	ng/ul	0.00

Target Compounds Qvalue

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

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