Quantitation Report (Qedit)

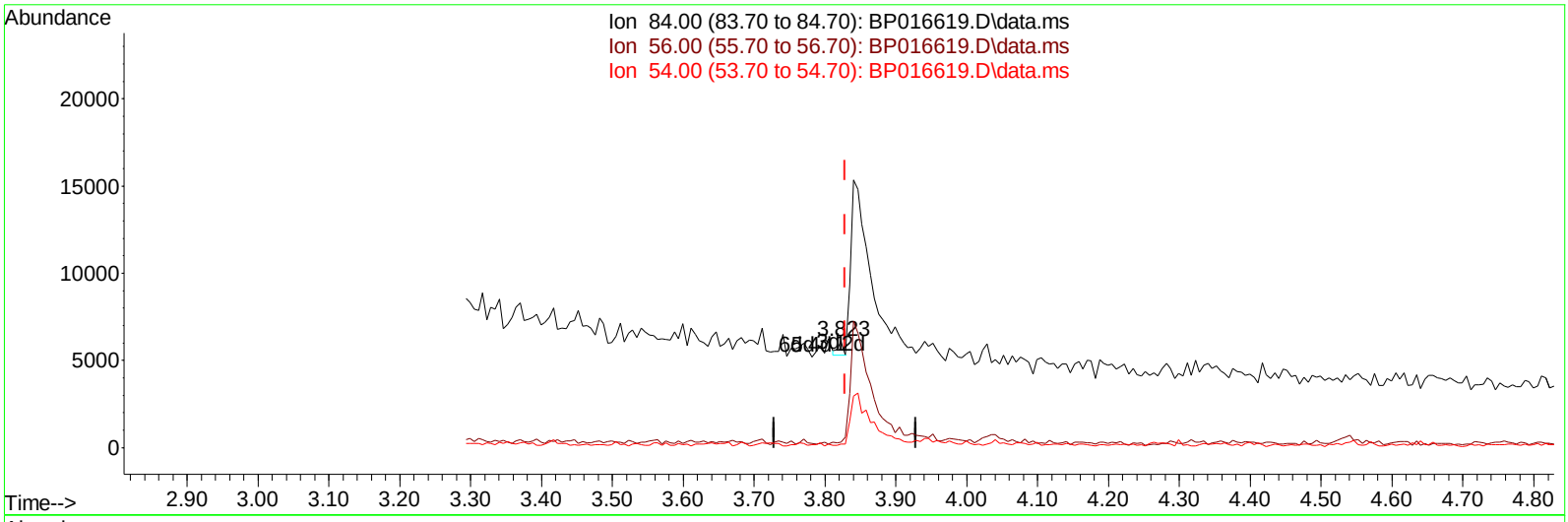
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
Data File : BP016619.D
Acq On : 17 Aug 2023 00:55
Operator : MA/JU
Sample : 03967-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

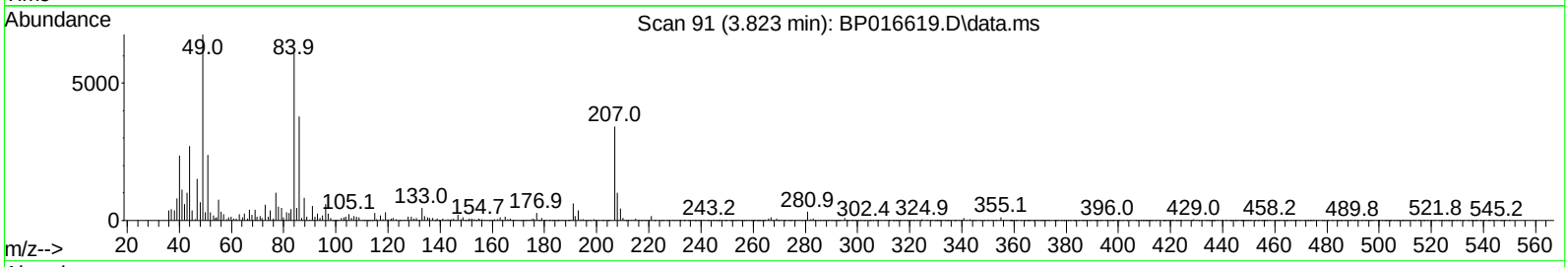
Quant Time: Aug 17 01:43:54 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Aug 17 01:38:34 2023
Response via : Initial Calibration

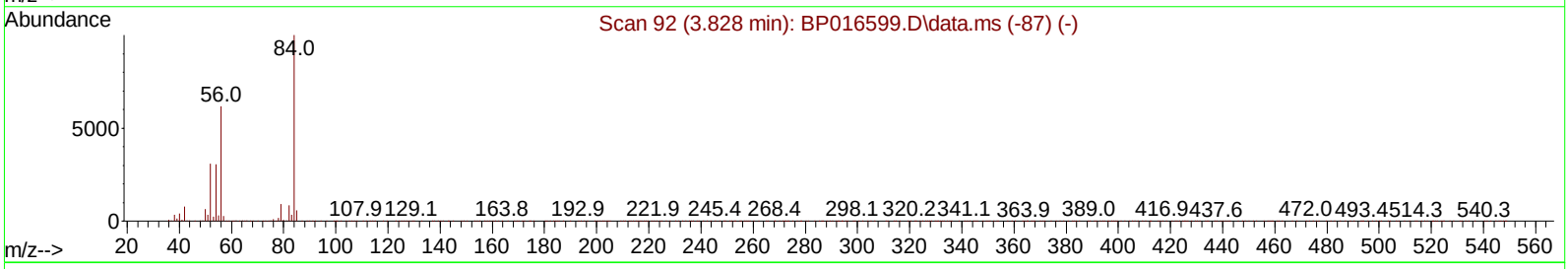
Instrument :
BNA_P
ClientSampleId :
A48M3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 08/17/2023
Supervised By :mohammad ahmed 08/17/2023







TIC: BP016619.D\data.ms

(4) Pyridine-d5 (S)

3.823min (-0.006) 0.01 ng/ul

response 402

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.90	5.37#
54.00	30.30	3.53#
0.00	0.00	0.00

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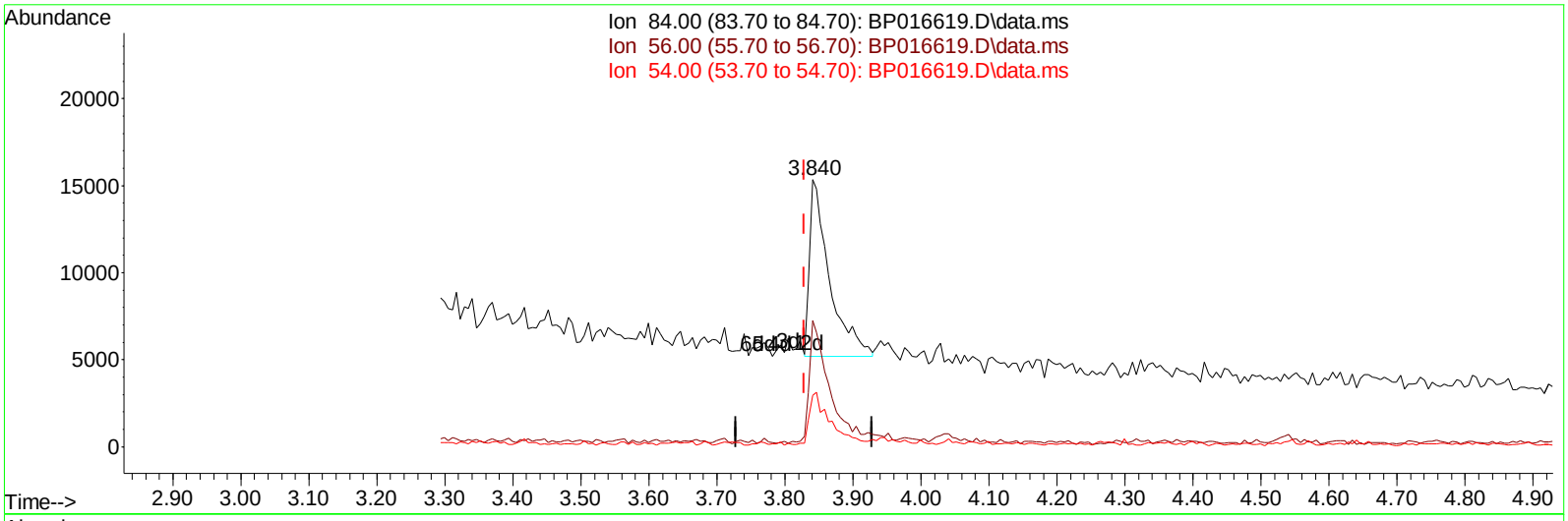

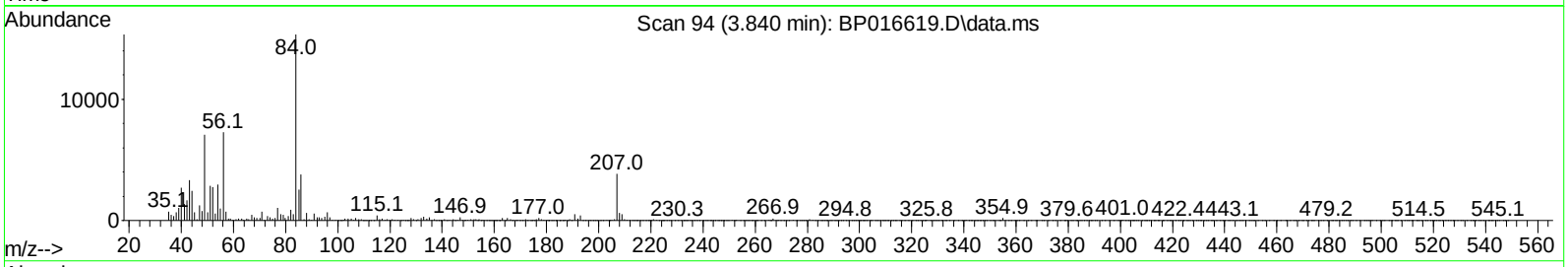

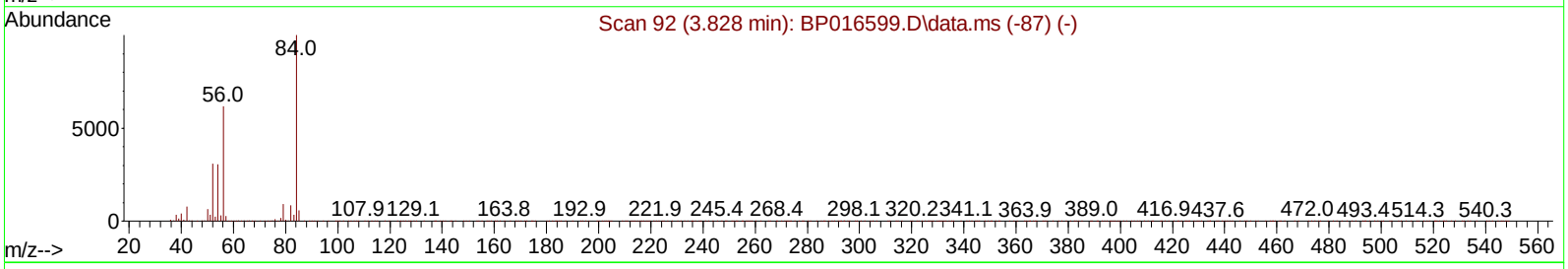

TIC: BP016619.D\data.ms

(4) Pyridine-d5 (S)

3.840min (+ 0.012) 0.76 ng/ul m

response 20704

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	61.90	47.40#
54.00	30.30	19.25#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP081623\
Data File : BP016619.D
Acq On : 17 Aug 2023 00:55
Operator : MA/JU
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Misc :
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Quant Time: Aug 17 01:45:25 2023
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Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.052	152	338774	20.000	ng/ul	0.00
20) Naphthalene-d8	10.887	136	1601341	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.704	164	1033465	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	2423733	20.000	ng/ul	0.00
79) Chrysene-d12	21.563	240	2515179	20.000	ng/ul	0.00
88) Perylene-d12	24.145	264	2813991	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	20673	2.319	ng/uL	0.00
4) Pyridine-d5	3.840	84	20704m	0.763	ng/ul	0.01
7) Phenol-d5	7.193	99	313559	10.040	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.387	67	219057	10.649	ng/ul	0.00
11) 2-Chlorophenol-d4	7.569	132	232145	10.348	ng/ul	0.00
15) 4-Methylphenol-d8	8.752	113	202036	8.135	ng/ul	0.00
21) Nitrobenzene-d5	9.252	128	120590	10.702	ng/ul	0.00
24) 2-Nitrophenol-d4	9.963	143	116234	10.861	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.493	165	220588	10.474	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	221958	6.142	ng/ul	0.00
46) Dimethylphthalate-d6	14.110	166	817043	10.930	ng/ul	0.00
49) Acenaphthylene-d8	14.404	160	914530	10.839	ng/ul	0.00
54) 4-Nitrophenol-d4	14.898	143	125816	8.262	ng/ul	0.00
60) Fluorene-d10	15.693	176	742463	11.651	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.816	200	71830	6.230	ng/ul	0.00
73) Anthracene-d10	17.569	188	1176186	11.219	ng/ul	0.00
81) Pyrene-d10	19.798	212	1574400	11.946	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.969	264	1578196	11.703	ng/ul	0.00
Target Compounds						Qvalue
16) Acetophenone	8.899	105	108953	2.717	ng/ul	98

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

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