Quantitation Report (Qedit)

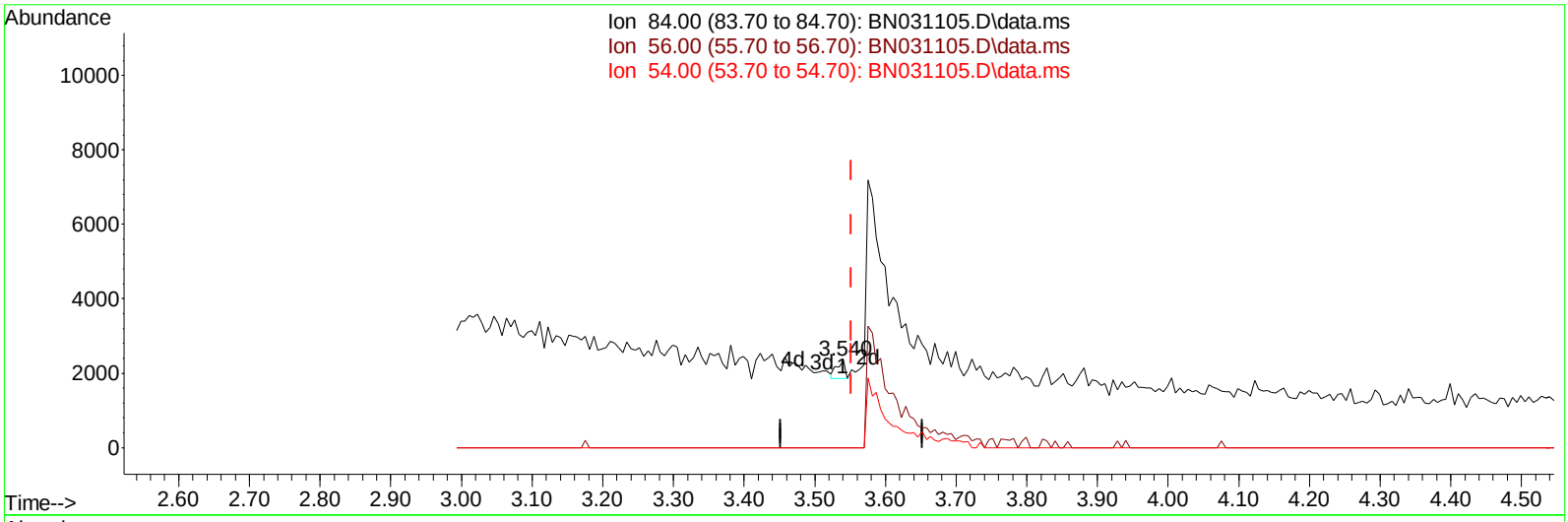
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031105.D
Acq On : 21 Apr 2024 12:30
Operator : MA/JU
Sample : P2161-01
Misc :
ALS Vial : 56 Sample Multiplier: 1

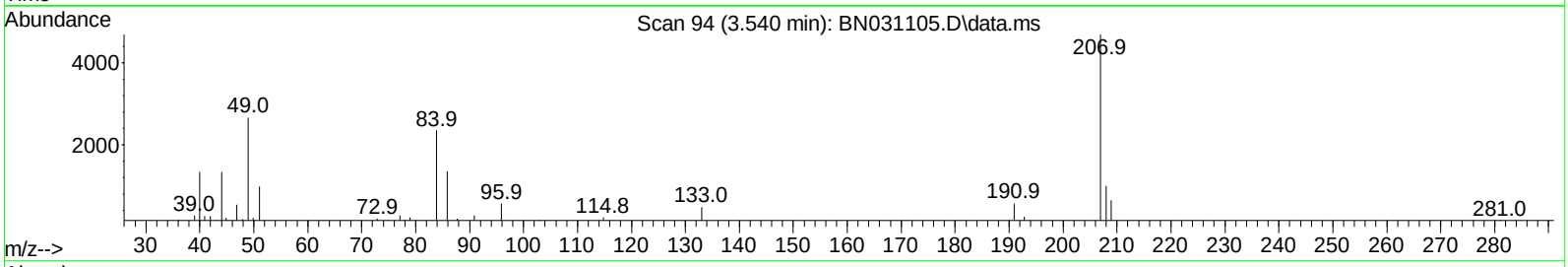
Quant Time: Apr 21 14:42:59 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 19 22:56:25 2024
Response via : Initial Calibration

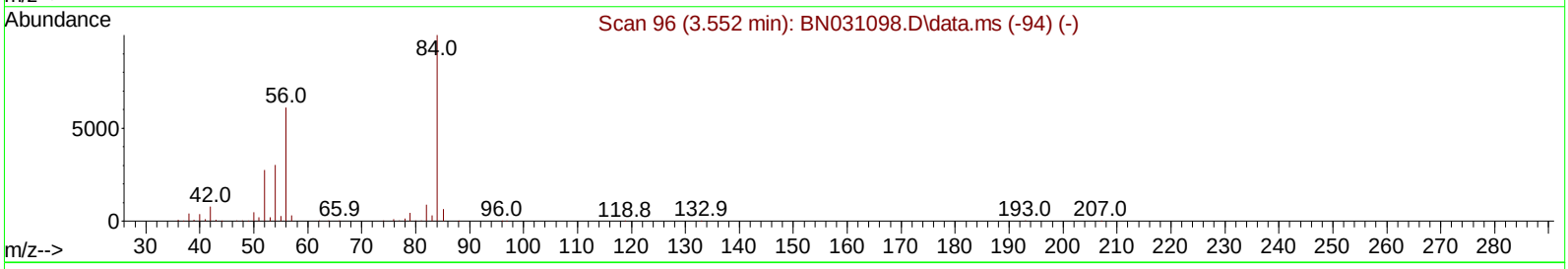
Instrument :
BNA_N
ClientSampleId :
C0RC3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/22/2024
Supervised By :mohammad ahmed 04/22/2024







TIC: BN031105.D\data.ms

(4) Pyridine-d5 (S)

3.540min (-0.012) 0.05 ng/ul

response 383

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.00	0.00#
54.00	30.10	0.00#
0.00	0.00	0.00

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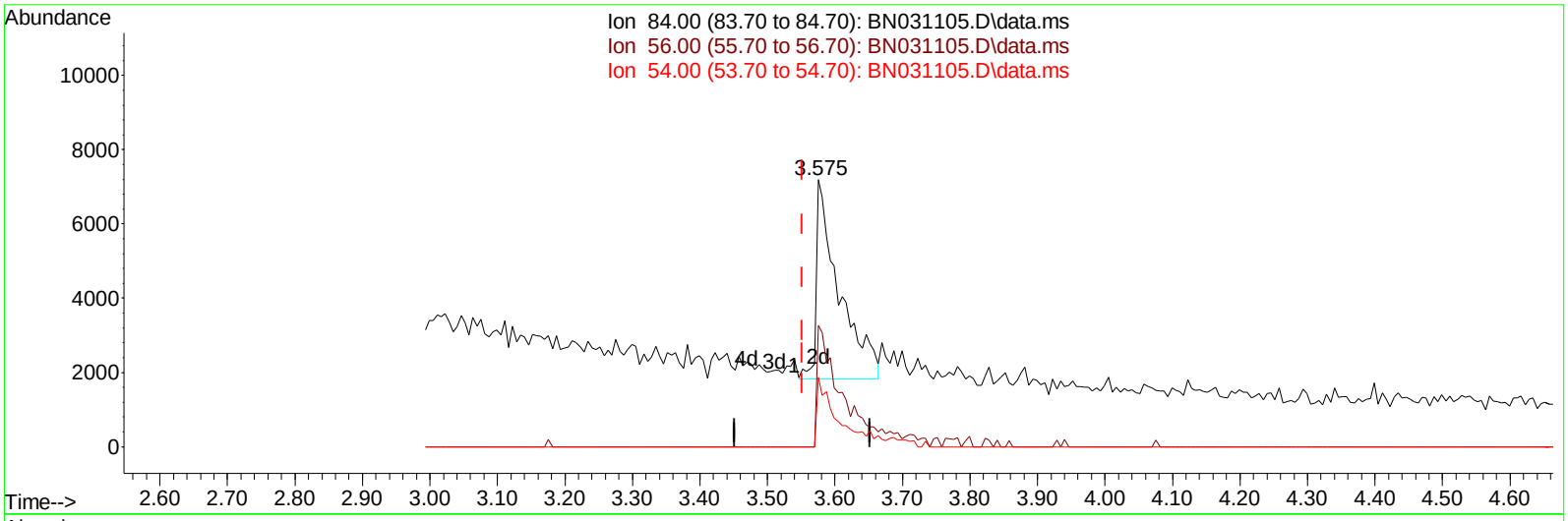

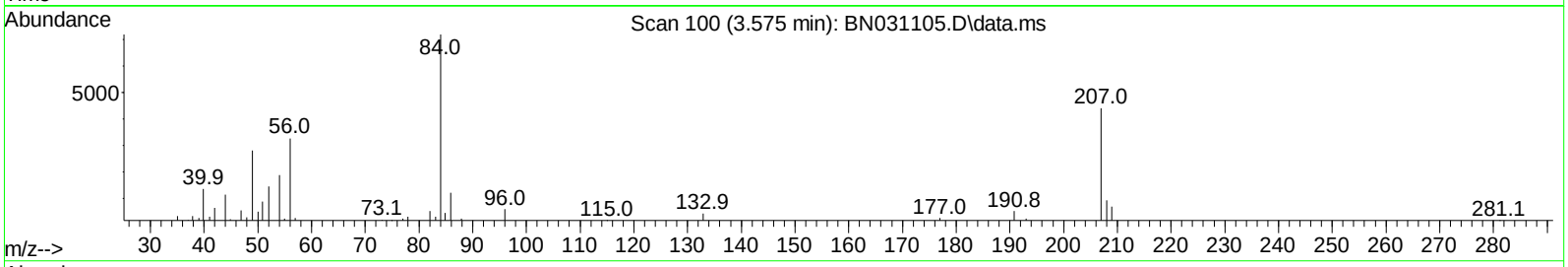

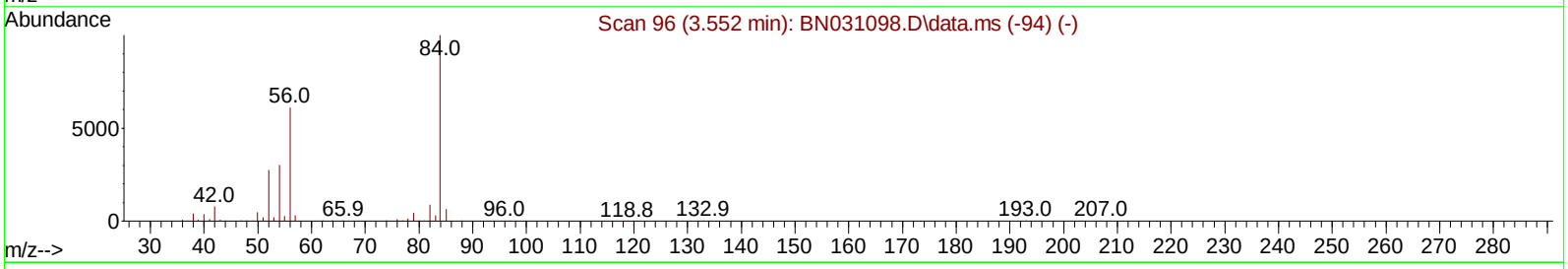

TIC: BN031105.D\data.ms

(4) Pyridine-d5 (S)

3.575min (+ 0.023) 1.65 ng/ul m

response 12591

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	59.00	45.33#
54.00	30.10	26.17
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Data File : BN031105.D
Acq On : 21 Apr 2024 12:30
Operator : MA/JU
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Misc :
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Quant Time: Apr 21 14:43:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 19 22:56:25 2024
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	124080	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.410	136	551629	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.281	164	378363	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.028	188	854006	20.000	ng/ul	0.00
79) Chrysene-d12	21.263	240	1059433	20.000	ng/ul	0.00
88) Perylene-d12	24.168	264	1444894	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	9369	3.528	ng/uL	0.00
4) Pyridine-d5	3.575	84	12591m	1.646	ng/ul	0.02
7) Phenol-d5	6.805	99	50872	4.860	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	172504	28.327	ng/ul	0.00
11) 2-Chlorophenol-d4	7.164	132	168416	20.665	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	105463	11.981	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128	113451	28.625	ng/ul	0.00
24) 2-Nitrophenol-d4	9.505	143	109515	26.582	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.034	165	197996	24.370	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	87899	6.923	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166	774917	29.432	ng/ul	0.00
49) Acenaphthylene-d8	13.969	160	882974	29.594	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	19834	3.609	ng/ul	0.00
60) Fluorene-d10	15.275	176	697818	30.445	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	90327	20.964	ng/ul	0.00
73) Anthracene-d10	17.128	188	1276009	34.472	ng/ul	0.00
81) Pyrene-d10	19.433	212	1707321	32.795	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.963	264	2308031	33.685	ng/ul	-0.01
Target Compounds						Qvalue

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

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