Quantitation Report (Qedit)

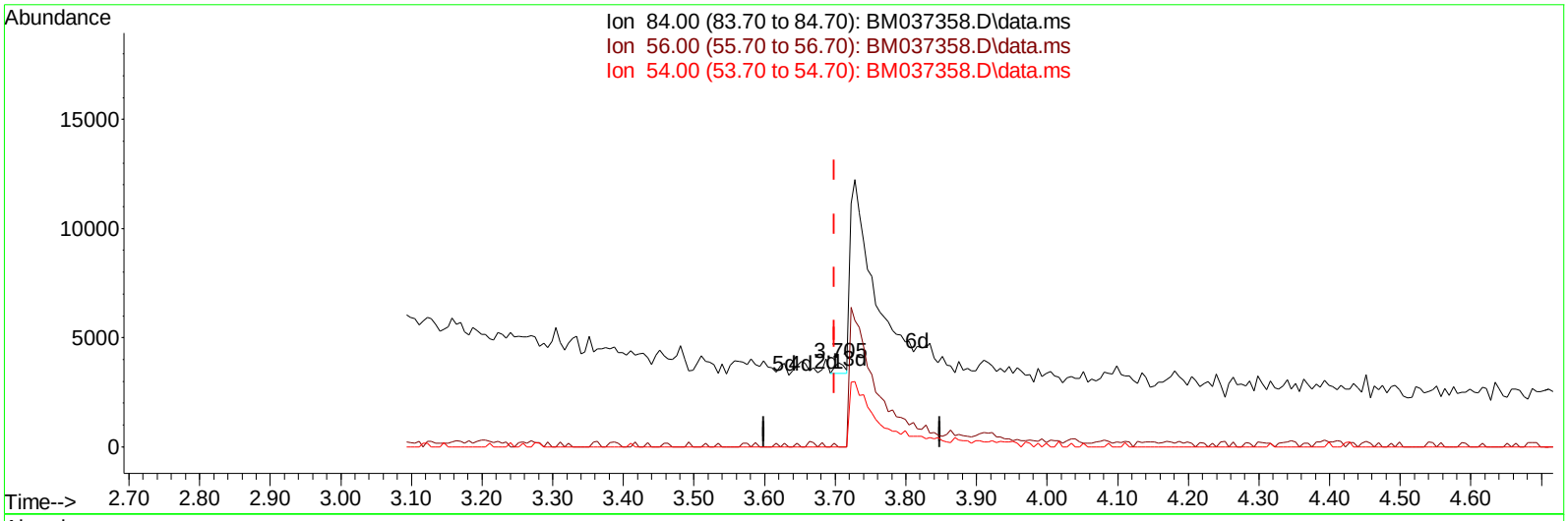
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
Data File : BM037358.D
Acq On : 03 Nov 2022 22:02
Operator : CG/JU
Sample : N5276-15
Misc :
ALS Vial : 21 Sample Multiplier: 1

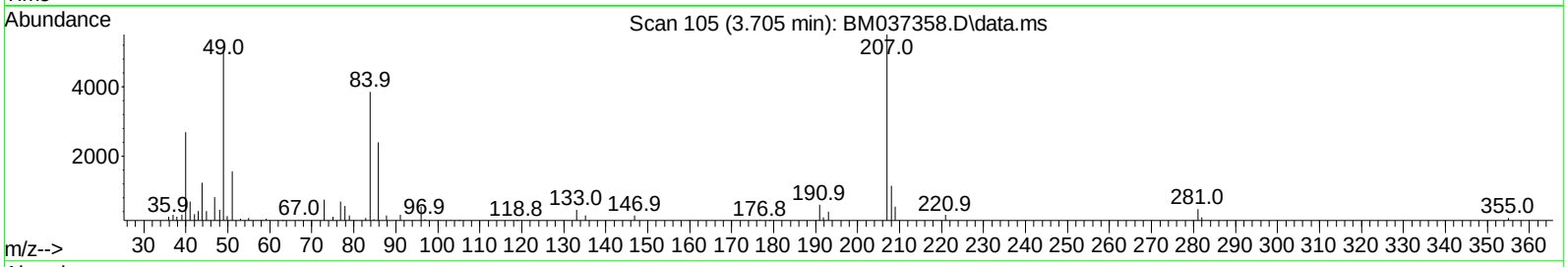
Quant Time: Nov 04 01:07:05 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 02 03:33:37 2022
Response via : Initial Calibration

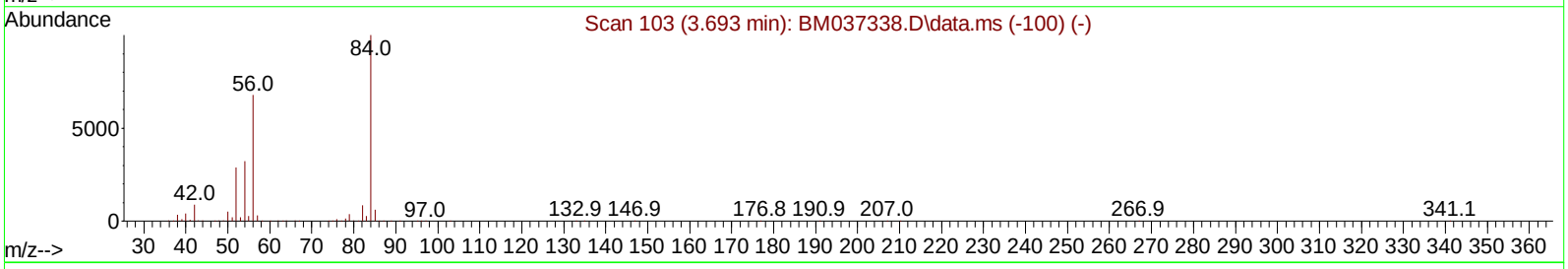
Instrument :
BNA_M
ClientSampleId :
EW4X5

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 11/04/2022
Supervised By :Jagrut Upadhyay 11/07/2022







TIC: BM037358.D\data.ms

(4) Pyridine-d5 (S)

3.705min (+ 0.006) 0.02 ng/ul

response 445

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	0.00#
54.00	30.90	0.00#
0.00	0.00	0.00

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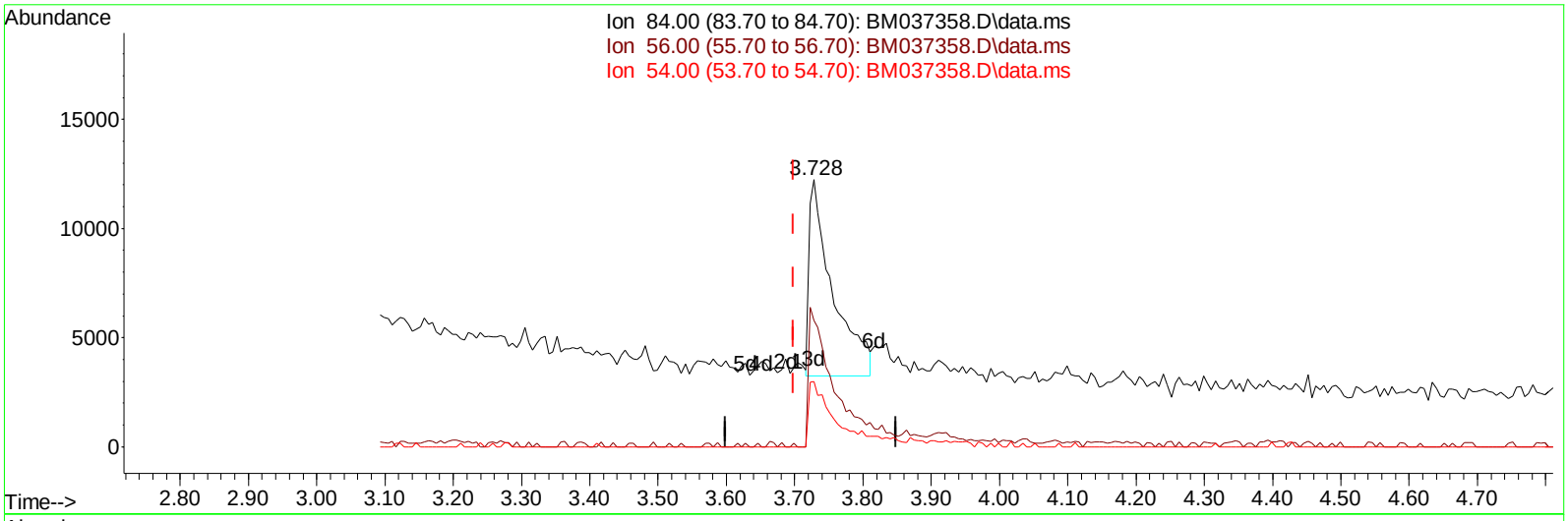

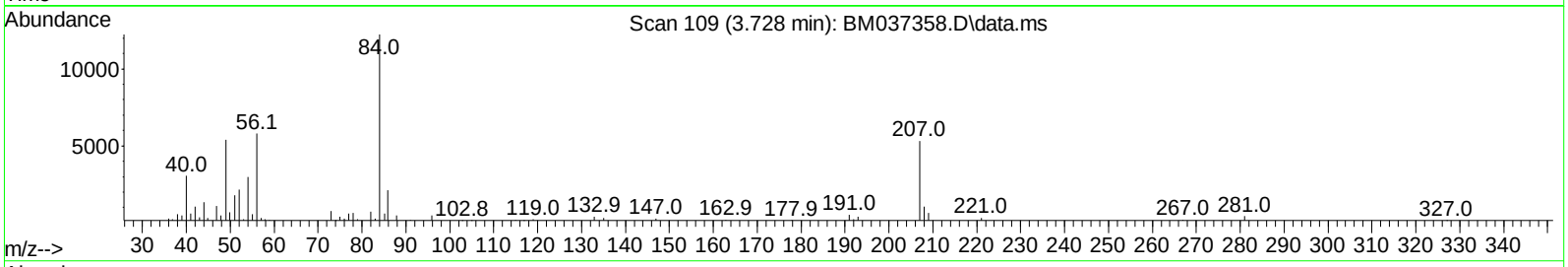

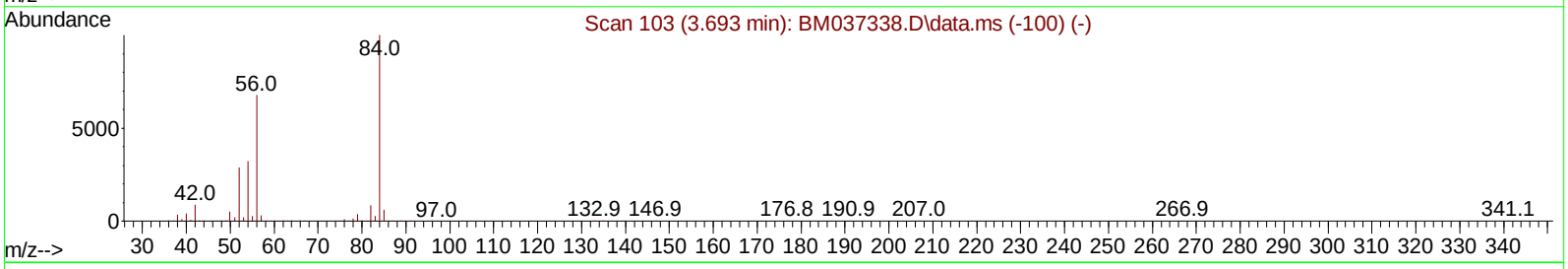

TIC: BM037358.D\data.ms

(4) Pyridine-d5 (S)

3.728min (+ 0.030) 1.03 ng/ul m

response 21680

Ion	Exp%	Act%
84.00	100.00	100.00
56.00	65.70	47.32#
54.00	30.90	24.44#
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.851	152	311126	20.000	ng/ul	0.00
20) Naphthalene-d8	10.651	136	1405299	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.486	164	917544	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.221	188	1897219	20.000	ng/ul	-0.01
79) Chrysene-d12	21.398	240	1712890	20.000	ng/ul	0.00
88) Perylene-d12	23.739	264	1481524	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	30907	4.379	ng/uL	0.00
4) Pyridine-d5	3.728	84	21680m	1.029	ng/ul	0.03
7) Phenol-d5	7.022	99	160734	5.936	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	475172	28.598	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	456813	22.081	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	299603	13.550	ng/ul	-0.01
21) Nitrobenzene-d5	9.016	128	303125	28.876	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.739	143	298911	26.141	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	506117	24.008	ng/ul	0.00
31) 4-Chloroaniline-d4	10.798	131	138291	4.293	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	1888048	29.761	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	2222454	29.139	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	50663	3.995	ng/ul	0.00
60) Fluorene-d10	15.474	176	1788987	31.622	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	267721	22.279	ng/ul	0.00
73) Anthracene-d10	17.321	188	2979071	34.095	ng/ul	-0.01
81) Pyrene-d10	19.609	212	3303225	36.904	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.592	264	2659385	34.872	ng/ul	-0.01

Target Compounds Qvalue

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

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