Quantitation Report (Qedit)

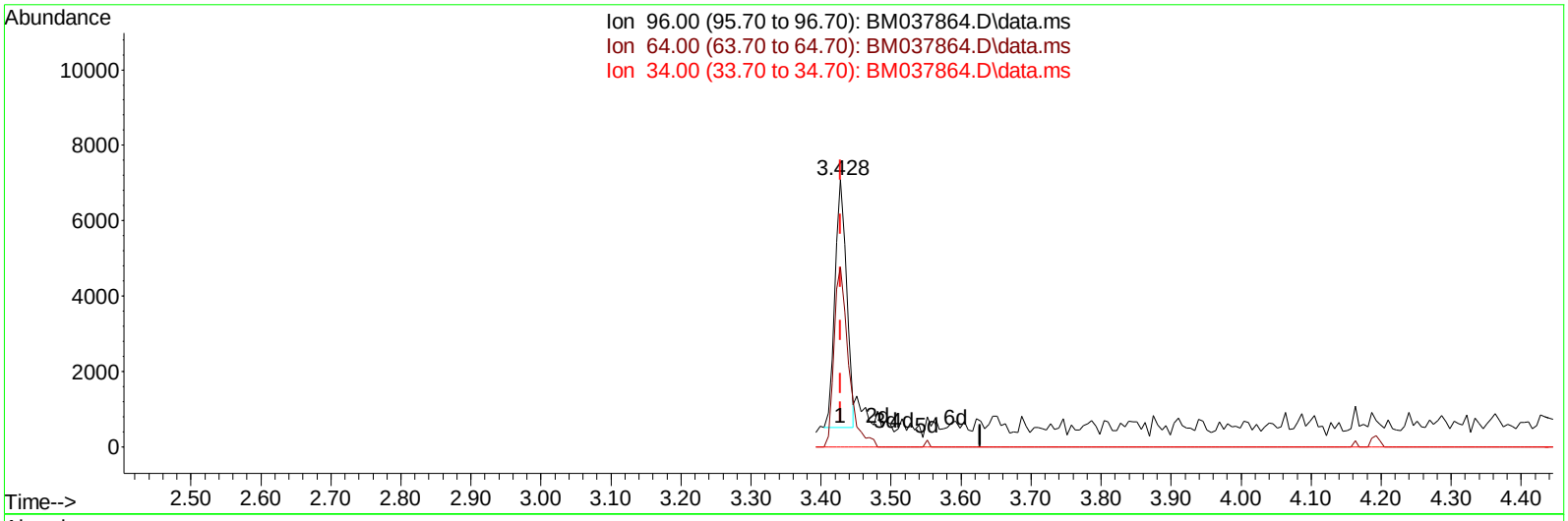
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037864.D
Acq On : 06 Dec 2022 21:12
Operator : CG/JU
Sample : N5738-12MEDL 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

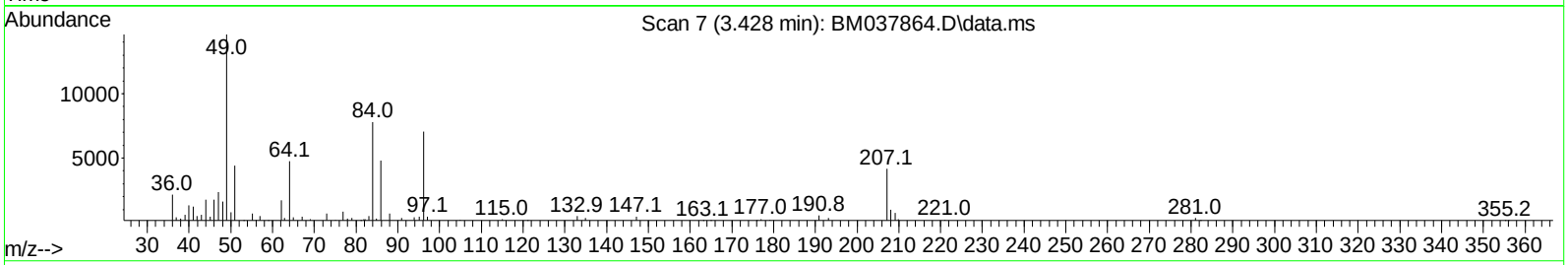
Quant Time: Dec 06 21:50:47 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Dec 02 01:14:59 2022
Response via : Initial Calibration

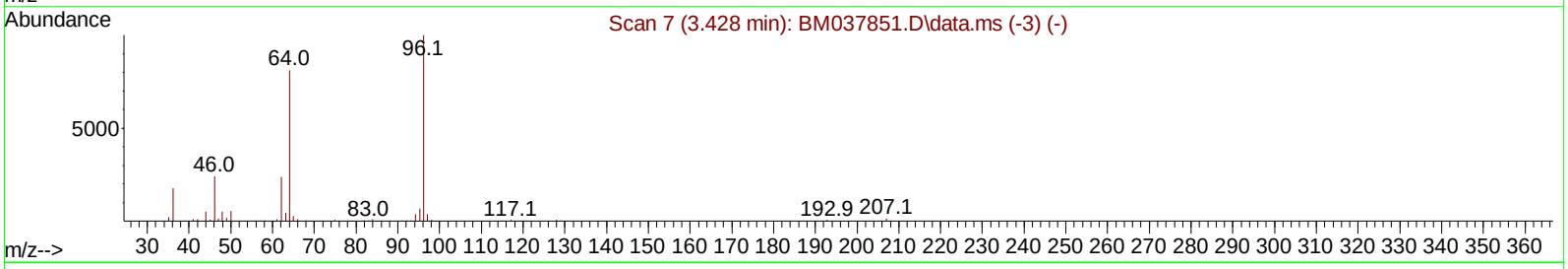
Instrument :
BNA_M
ClientSampleId :
GBNJ1MEDL

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 12/07/2022
Supervised By :mohammad ahmed 12/08/2022







TIC: BM037864.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.428min (-0.000) 1.35 ng/uL

response 7680

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	82.40	67.24
34.00	0.00	0.00
0.00	0.00	0.00

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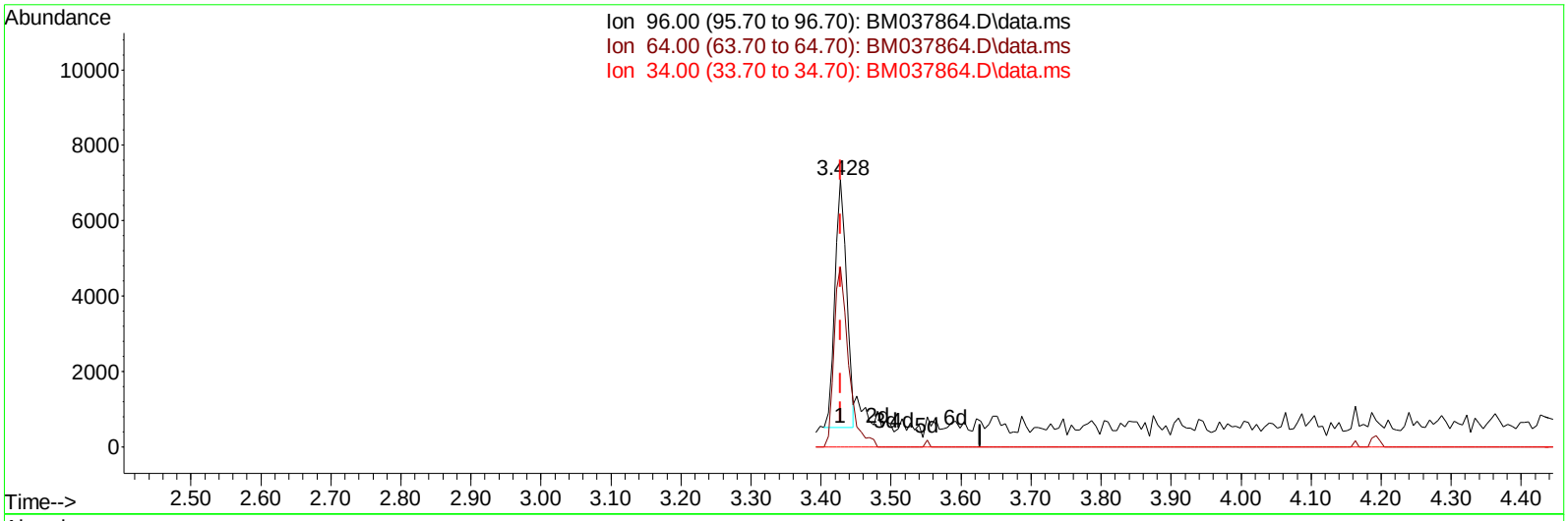

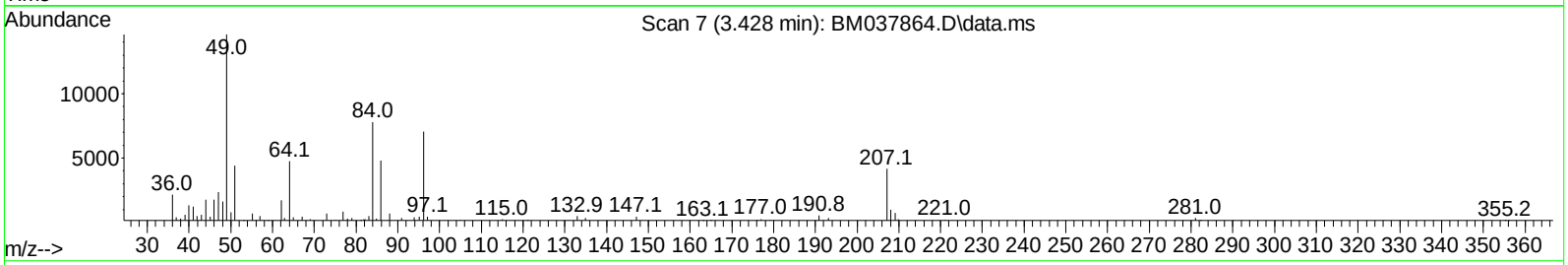

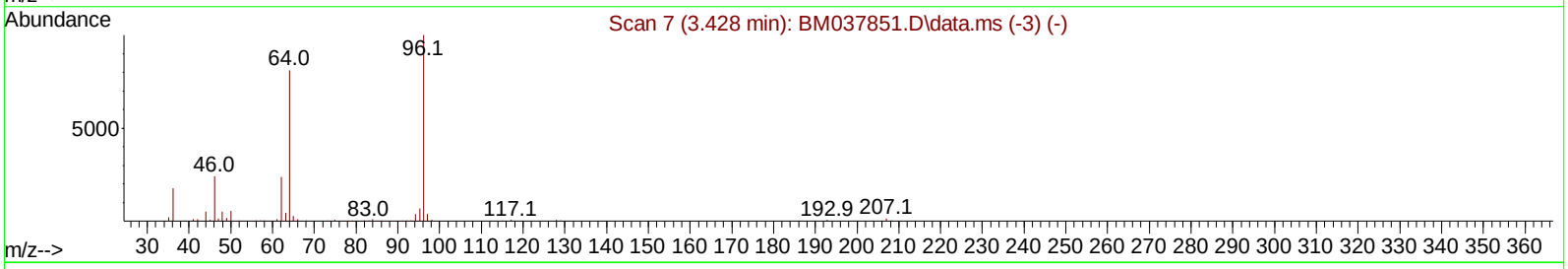

TIC: BM037864.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.428min (-0.000) 1.35 ng/uL

response 7680

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	82.40	67.24
34.00	0.00	0.00
0.00	0.00	0.00

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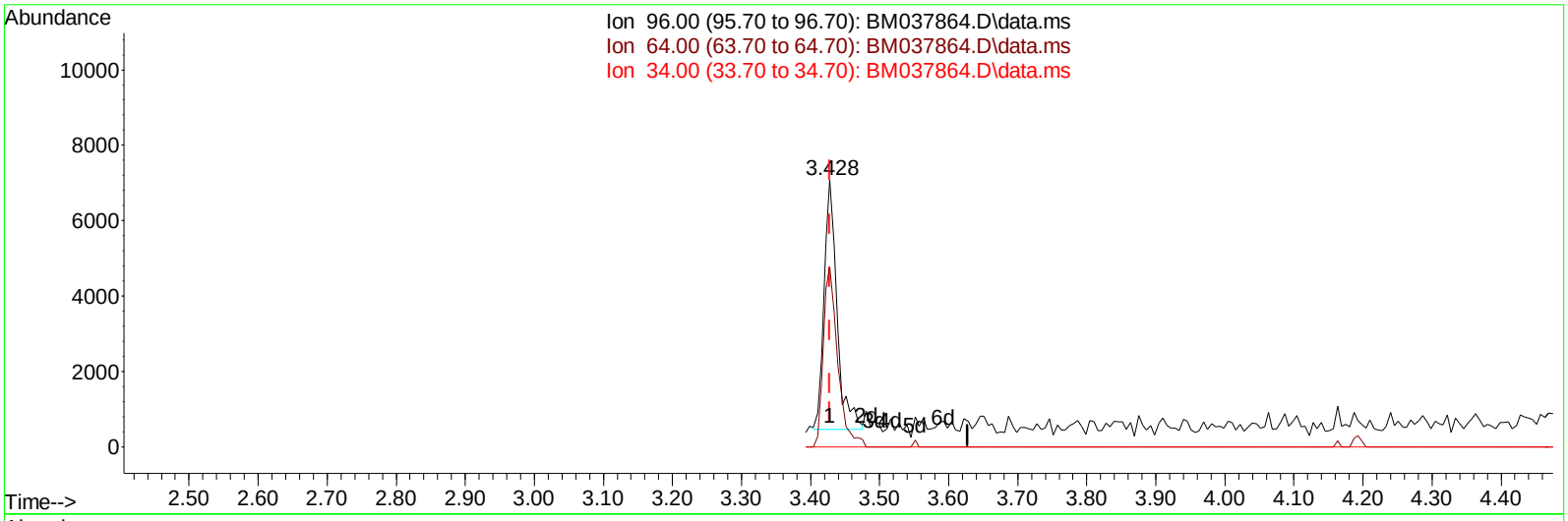

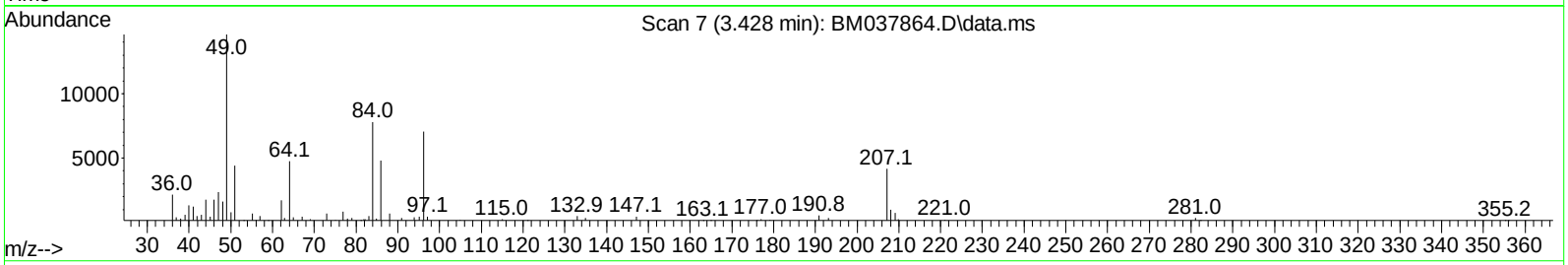

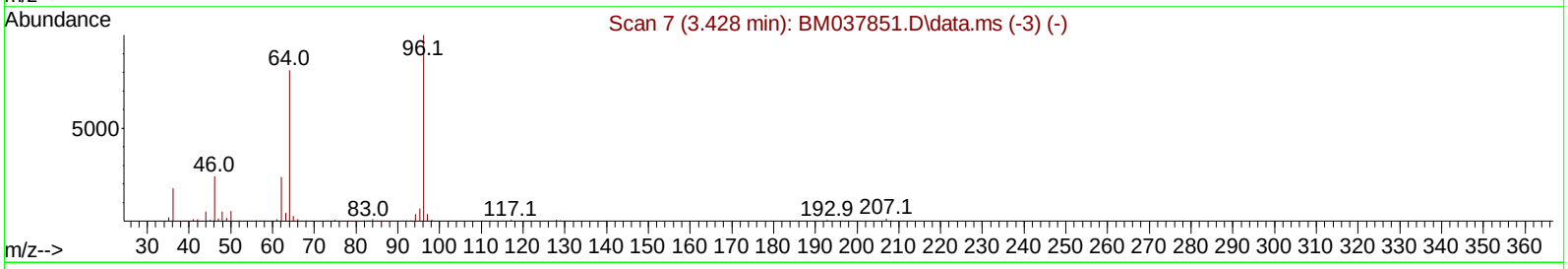

TIC: BM037864.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.428min (-0.000) 1.52 ng/uL m

response 8598

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	82.40	67.24
34.00	0.00	0.00
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	8.092	152	266013	20.000	ng/ul	0.00
20) Naphthalene-d8	10.910	136	1037675	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.721	164	573841	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.456	188	1243124	20.000	ng/ul	0.00
79) Chrysene-d12	21.621	240	1107664	20.000	ng/ul	-0.01
88) Perylene-d12	24.109	264	1223851	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.428	96	8598m	1.516	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.239	99	199785	8.336	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.410	67	132874	9.503	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.616	132	170918	9.432	ng/ul	0.00
15) 4-Methylphenol-d8	8.798	113	143866	7.548	ng/ul	0.00
21) Nitrobenzene-d5	9.263	128	83711	10.156	ng/ul	0.00
24) 2-Nitrophenol-d4	9.986	143	79752	9.350	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.527	165	145306	9.186	ng/ul	0.00
31) 4-Chloroaniline-d4	11.051	131	67194	2.769	ng/ul	0.00
46) Dimethylphthalate-d6	14.121	166	387674	9.838	ng/ul	-0.01
49) Acenaphthylene-d8	14.415	160	501364	10.462	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	50216	6.140	ng/ul	0.00
60) Fluorene-d10	15.704	176	407283	11.393	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.821	200	42931	6.556	ng/ul	0.00
73) Anthracene-d10	17.556	188	683529	11.956	ng/ul	0.00
81) Pyrene-d10	19.833	212	784461	13.672	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.938	264	726875	11.855	ng/ul	-0.02
Target Compounds						Qvalue
71) Pentachlorophenol	17.109	266	855085	93.327	ng/ul	99

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

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