Quantitation Report (Qedit)

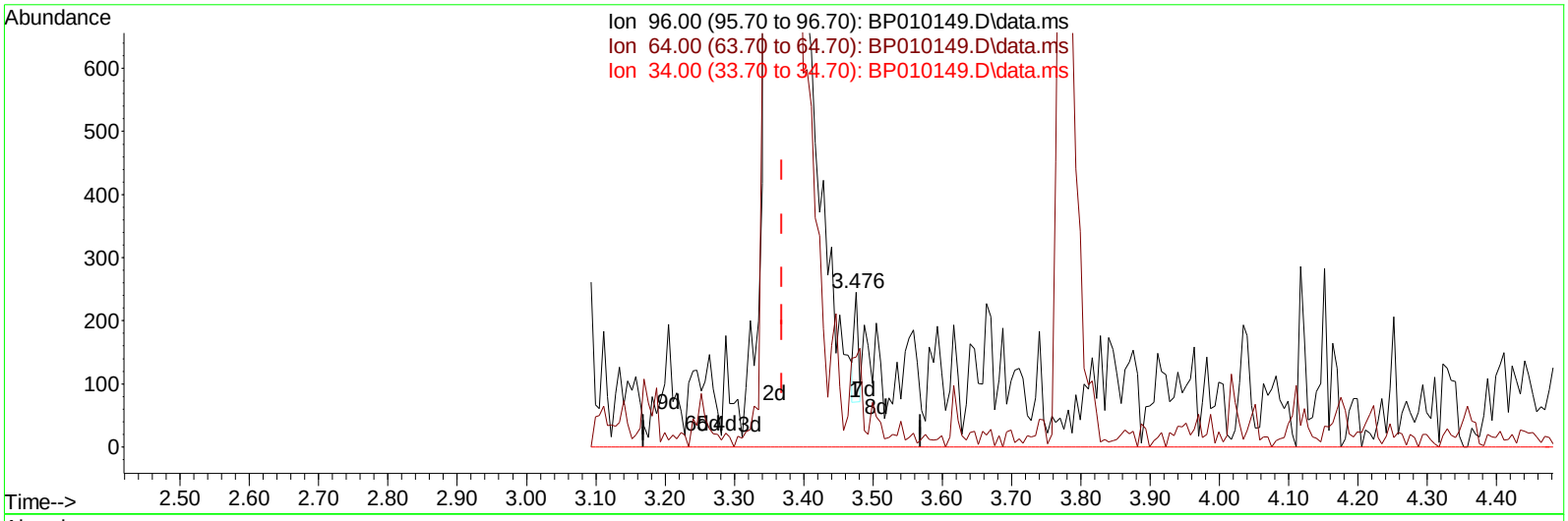
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042922\
Data File : BP010149.D
Acq On : 29 Apr 2022 13:55
Operator : CG/JU
Sample : SP5886
Misc : SFAM SURROGATES
ALS Vial : 9 Sample Multiplier: 1

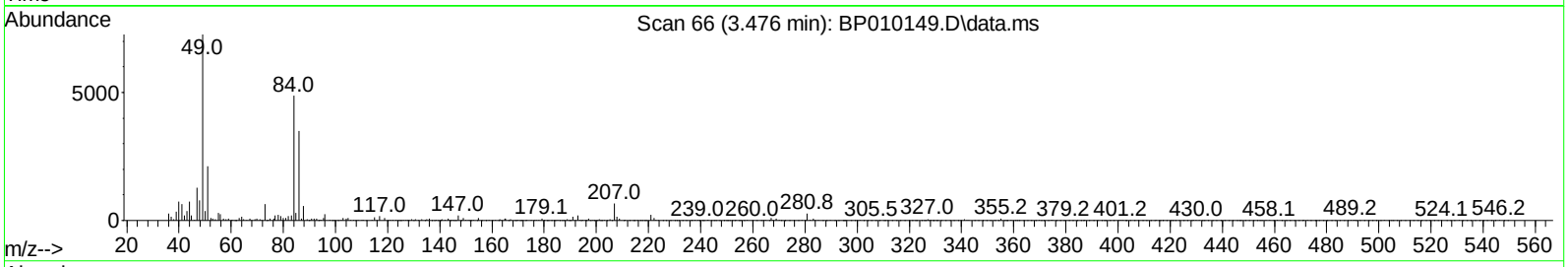
Quant Time: Apr 30 00:28:08 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration

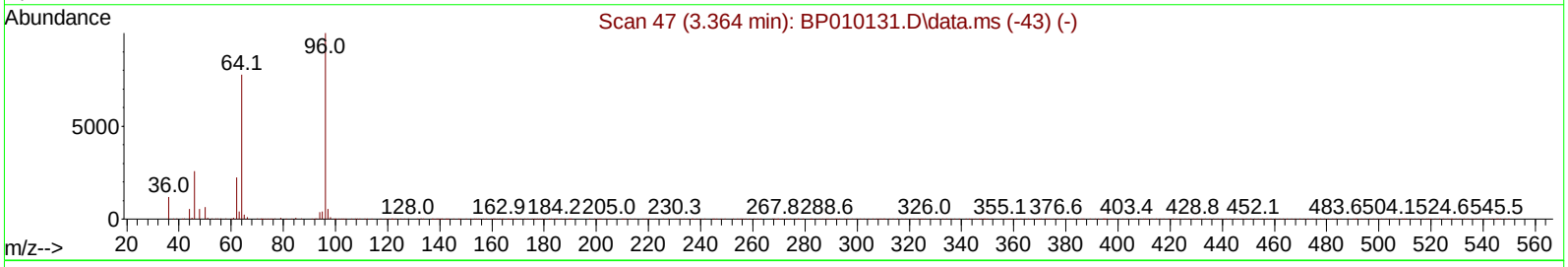
Instrument :
BNA_P
ClientSampleId :
SP5886

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/02/2022
Supervised By :Yogesh Patel 05/05/2022







TIC: BP010149.D\data.ms

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

3.476min (+ 0.107) 0.02 ng/uL

response 71

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	58.37
34.00	0.00	0.00
0.00	0.00	0.00

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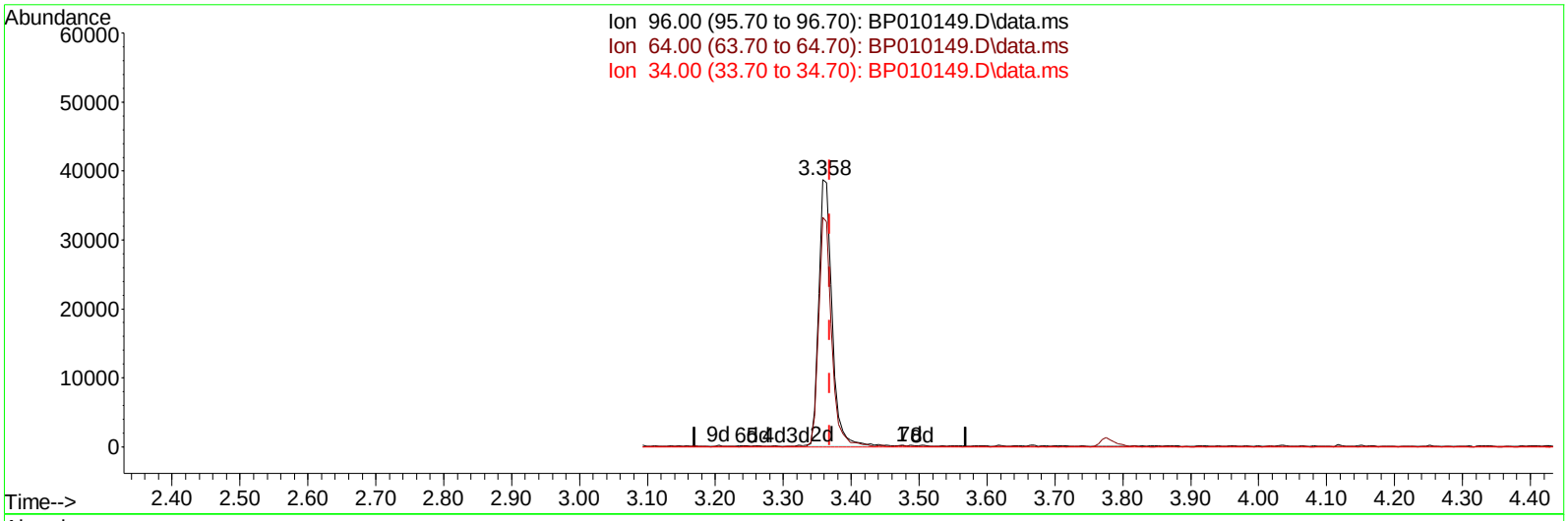

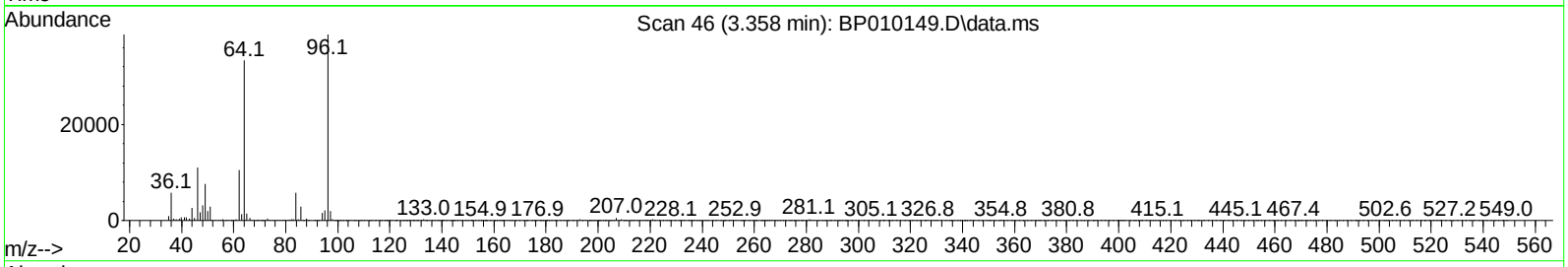

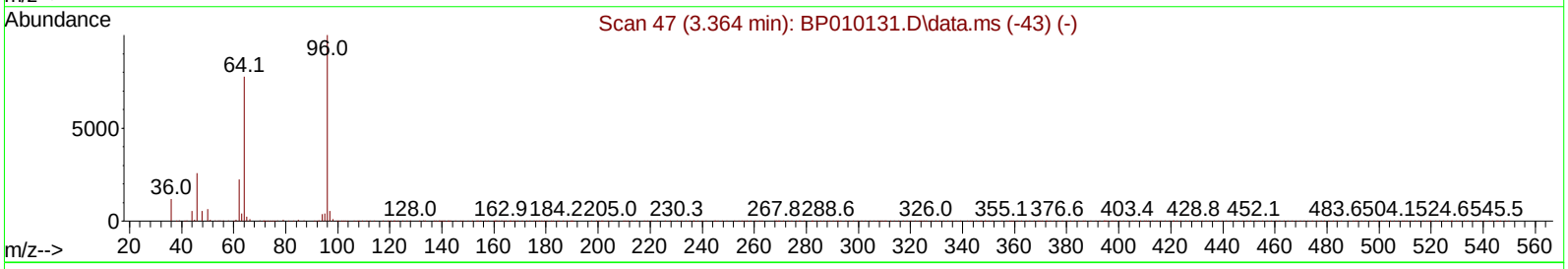

TIC: BP010149.D\data.ms

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

3.358min (-0.011) 17.19 ng/uL m

response 52549

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	86.02#
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	7.928	152	133273	20.000	ng/ul	#-0.01
20) Naphthalene-d8	10.728	136	525310	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.557	164	317865	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.310	188	631569	20.000	ng/ul	#-0.01
79) Chrysene-d12	21.398	240	605187	20.000	ng/ul	#-0.01
88) Perylene-d12	23.845	264	616043	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.358	96	52549m	17.194	ng/uL	-0.01
4) Pyridine-d5	3.776	84	728223	84.699	ng/ul	-0.01
7) Phenol-d5	7.093	99	875629	72.232	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.252	67	581695	79.969	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.458	132	740456	82.477	ng/ul	-0.01
15) 4-Methylphenol-d8	8.634	113	715600	71.248	ng/ul	-0.02
21) Nitrobenzene-d5	9.087	128	366695	87.767	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.811	143	412104	90.315	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.352	165	666789	75.064	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.863	131	934332	73.357	ng/ul	-0.02
46) Dimethylphthalate-d6	13.969	166	1969346	79.646	ng/ul	-0.01
49) Acenaphthylene-d8	14.251	160	2355266	78.704	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.757	143	311332	68.394	ng/ul	0.00
60) Fluorene-d10	15.557	176	1673819	78.695	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.669	200	322524	80.762	ng/ul	-0.01
73) Anthracene-d10	17.410	188	2479325	81.144	ng/ul	-0.01
81) Pyrene-d10	19.645	212	2789777	81.058	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	2651968	82.113	ng/ul	0.00
Target Compounds						Qvalue

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

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