Quantitation Report (Qedit)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011336.D
Acq On : 09 Aug 2022 18:59
Operator : CG/JU
Sample : N3900-16
Misc :
ALS Vial : 17 Sample Multiplier: 1

Quant Time: Aug 09 23:14:55 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

Instrument :
BNA_P
ClientSampleId :
C0CB3

Manual IntegrationsAPPROVED

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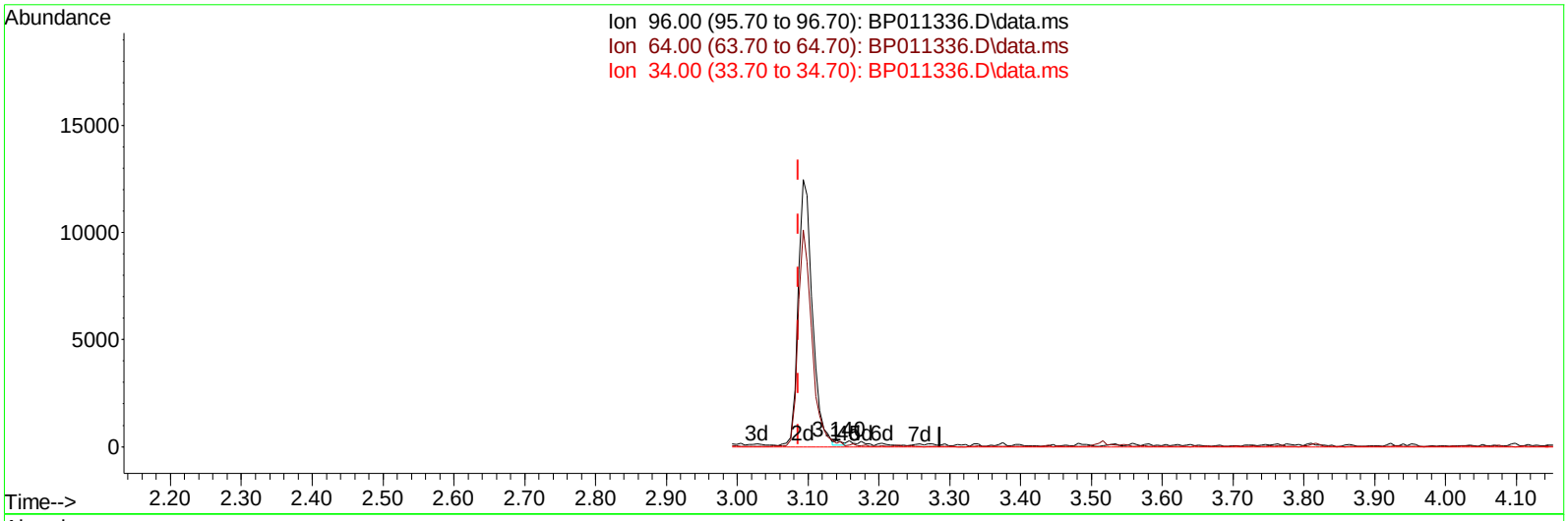
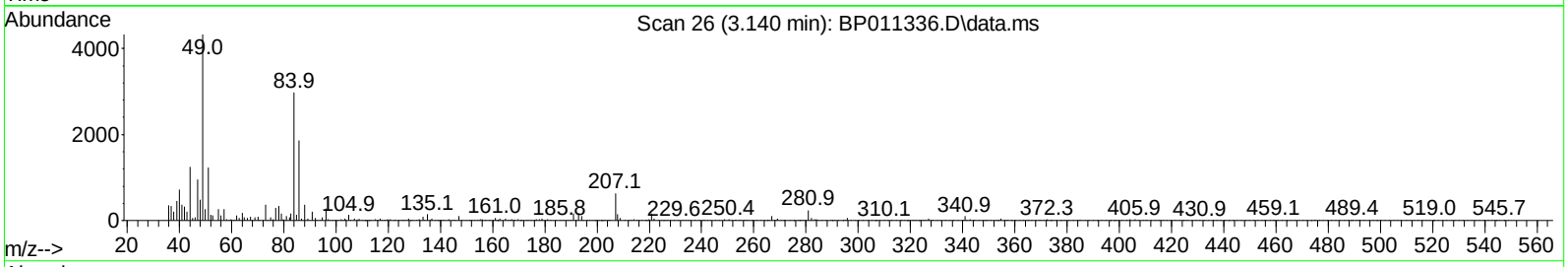
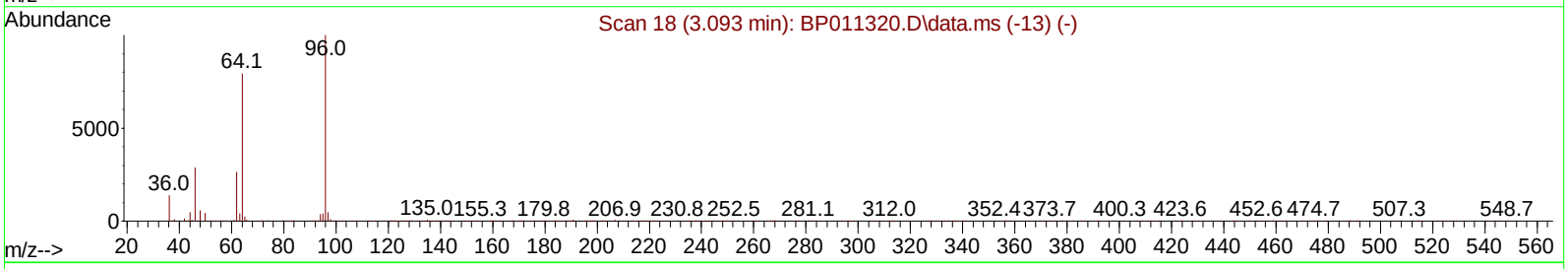
Supervised By :Sohil Jodhani 08/10/2022







TIC: BP011336.D\data.ms

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

3.140min (+ 0.053) 0.04 ng/uL

response 155

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	63.64
34.00	0.00	0.00
0.00	0.00	0.00

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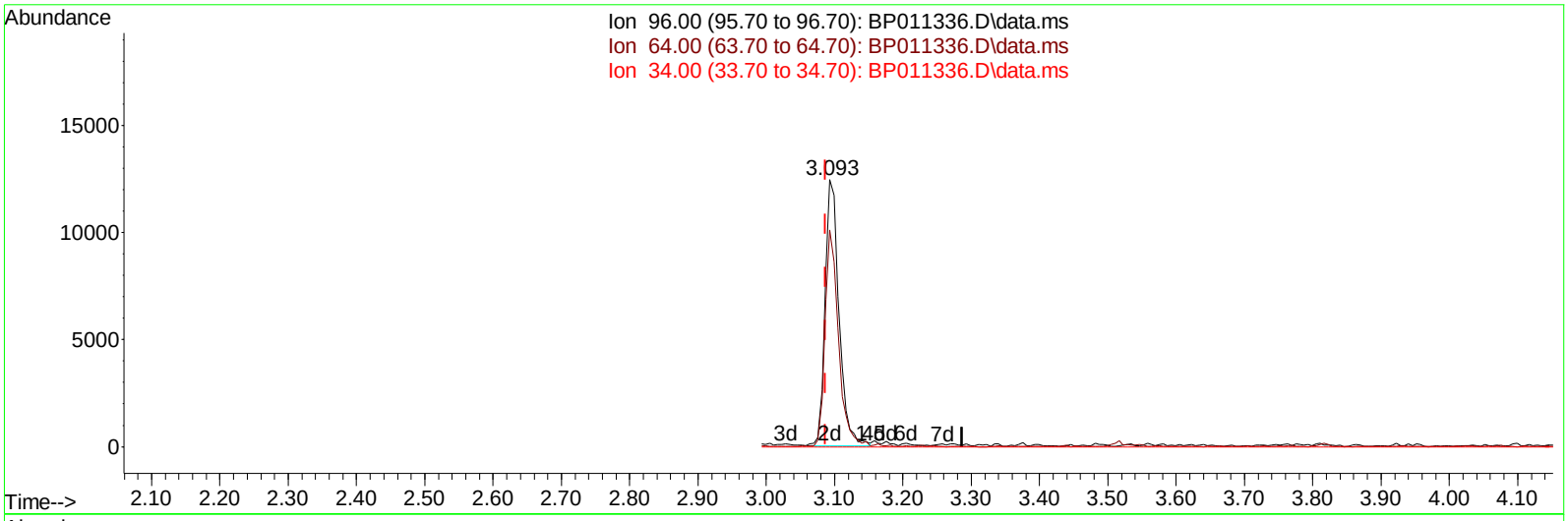

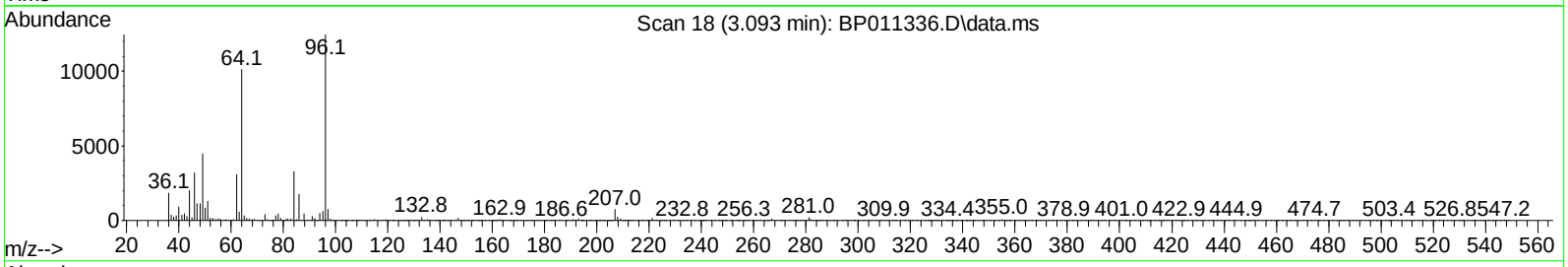

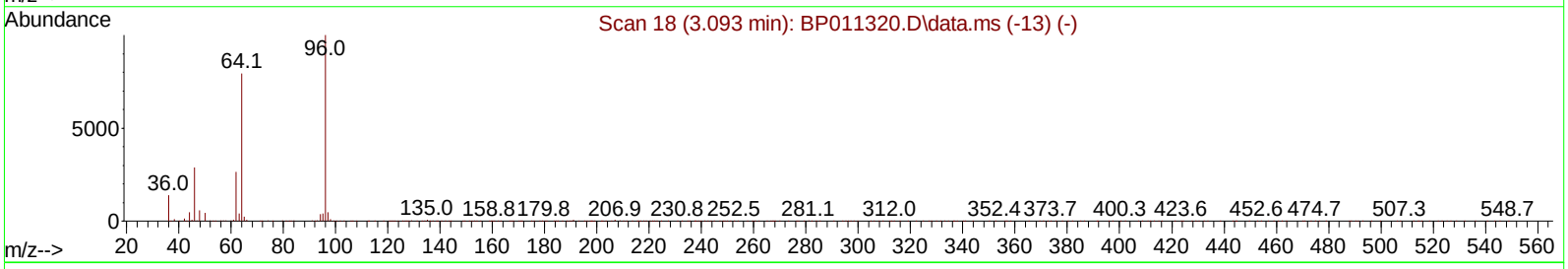

TIC: BP011336.D\data.ms

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

3.093min (+ 0.006) 4.07 ng/uL m

response 17653

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	81.33#
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	7.593	152	132996	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	561084	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	327543	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.034	188	655858	20.000	ng/ul	0.00
79) Chrysene-d12	21.157	240	607901	20.000	ng/ul	-0.01
88) Perylene-d12	23.468	264	615849	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	17653m	4.069	ng/uL	0.00
4) Pyridine-d5	3.511	84	77780	7.107	ng/ul	0.01
7) Phenol-d5	6.775	99	67245	5.713	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	243845	29.219	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	210444	22.788	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	134657	14.571	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	136352	35.707	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	125505	35.899	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	215266	30.490	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	320143	25.937	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	902207	39.698	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	1007083	36.812	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	14864	3.581	ng/ul	0.00
60) Fluorene-d10	15.263	176	742036	38.179	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	98890	32.462	ng/ul	0.00
73) Anthracene-d10	17.134	188	1164270	39.494	ng/ul	0.00
81) Pyrene-d10	19.392	212	1327742	39.644	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	1224352	39.422	ng/ul	0.00
Target Compounds						Qvalue
6) Benzaldehyde	6.746	77	26852	5.235	ng/ul	97
16) Acetophenone	8.422	105	27047	1.809	ng/ul	97
35) 4-Chloro-3-methylphenol	11.698	107	15268	1.759	ng/ul	98
59) Diethylphthalate	15.110	149	61098	2.562	ng/ul	98

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

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