Quantitation Report (Qedit)

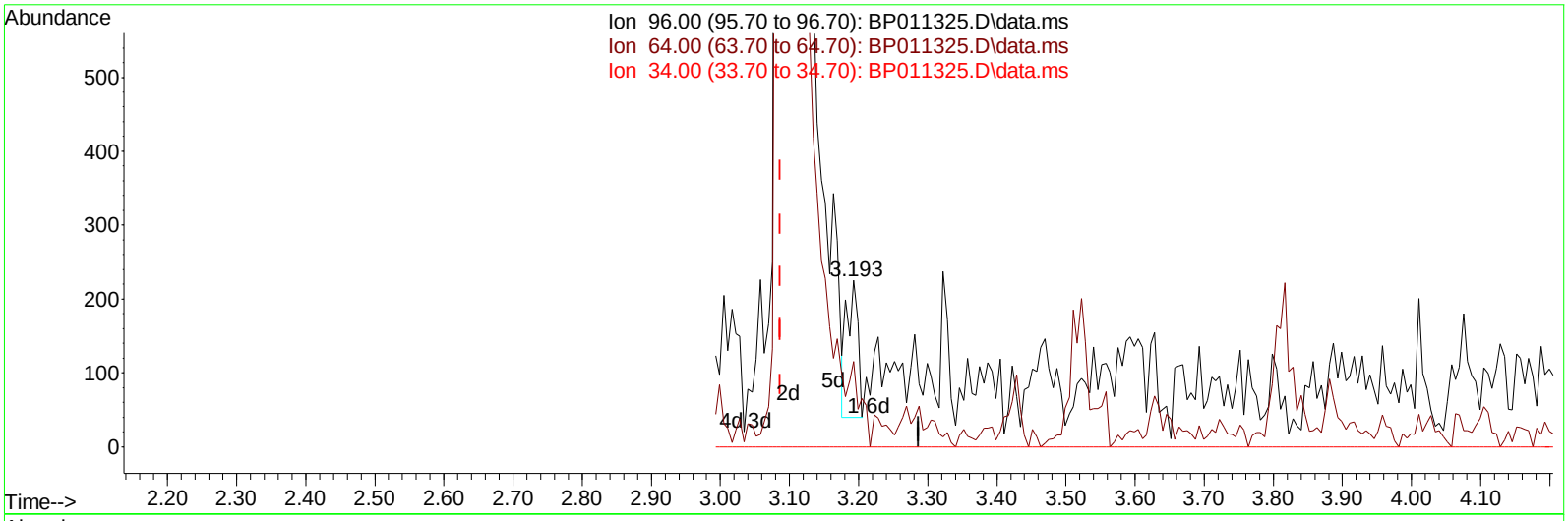
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080922\
Data File : BP011325.D
Acq On : 09 Aug 2022 12:35
Operator : CG/JU
Sample : N4026-19
Misc :
ALS Vial : 7 Sample Multiplier: 1

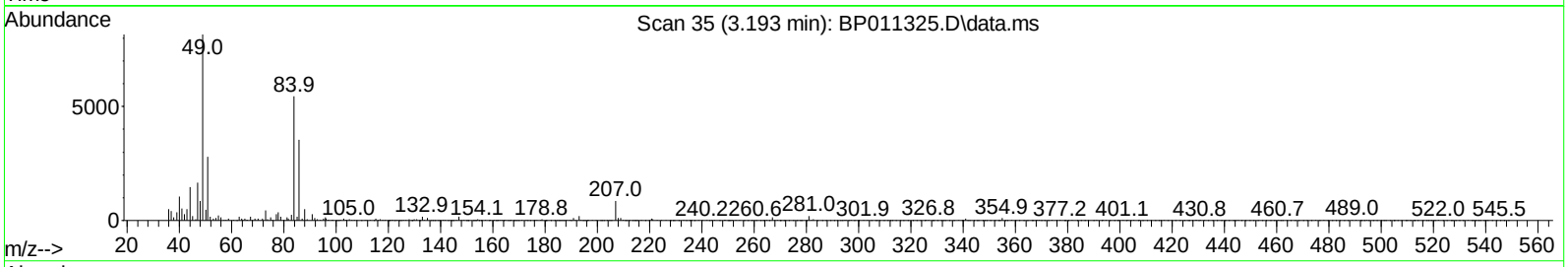
Quant Time: Aug 09 23:10:10 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080822.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 18:47:19 2022
Response via : Initial Calibration

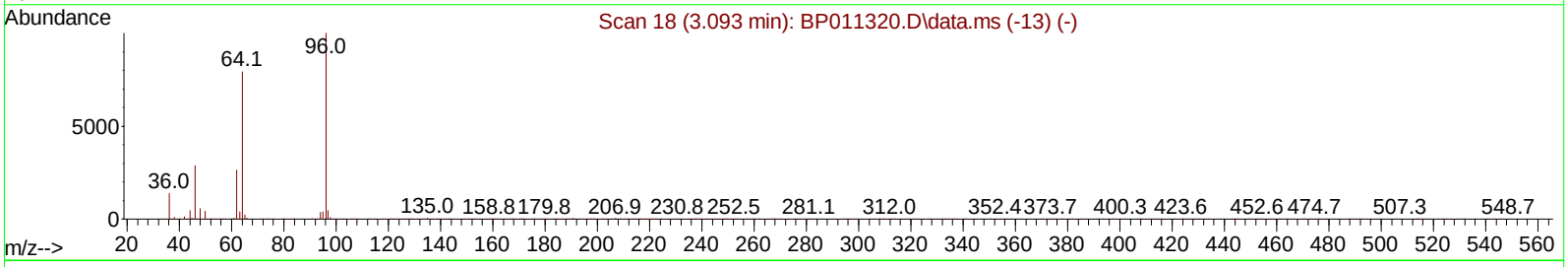
Instrument :
BNA_P
ClientSampleId :
GBJJ3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/10/2022
Supervised By :Sohil Jodhani 08/10/2022







TIC: BP011325.D\data.ms

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

3.193min (+ 0.106) 0.04 ng/uL

response 206

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	51.56
34.00	0.00	0.00
0.00	0.00	0.00

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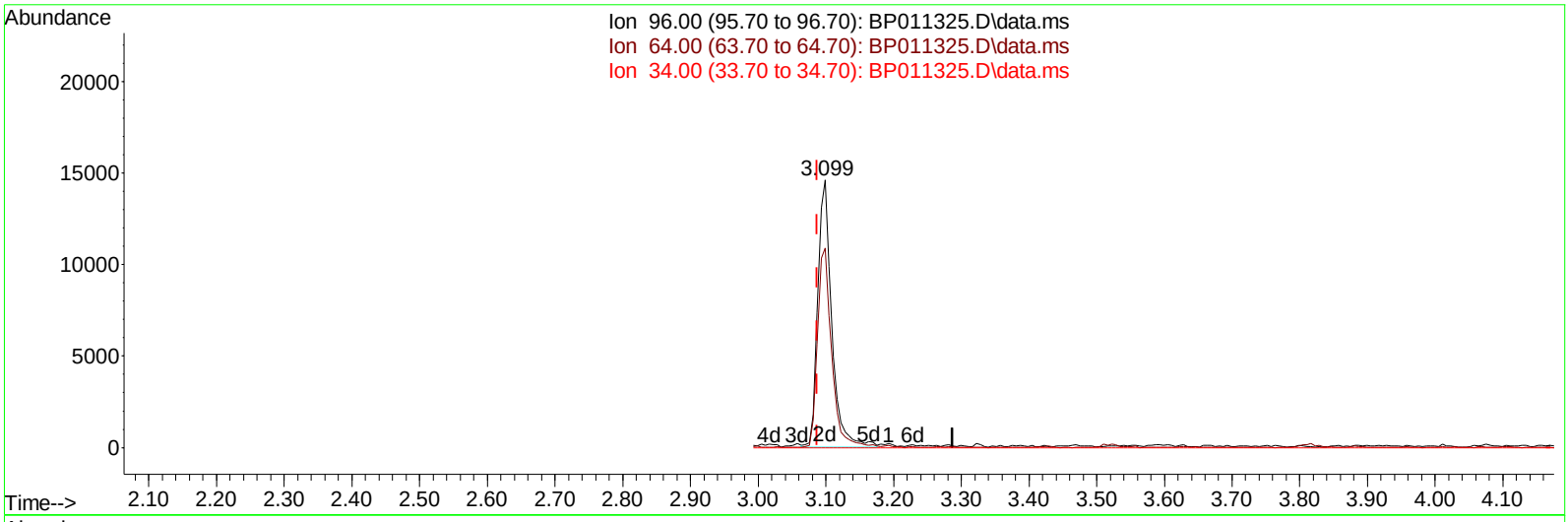

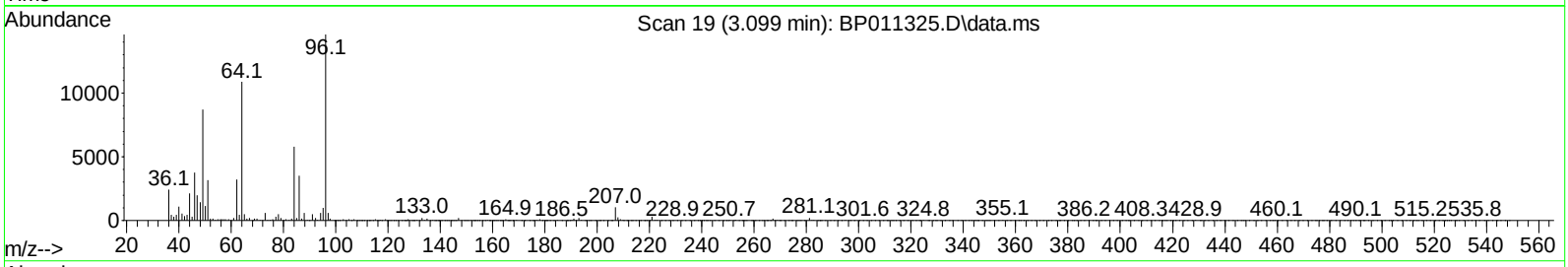

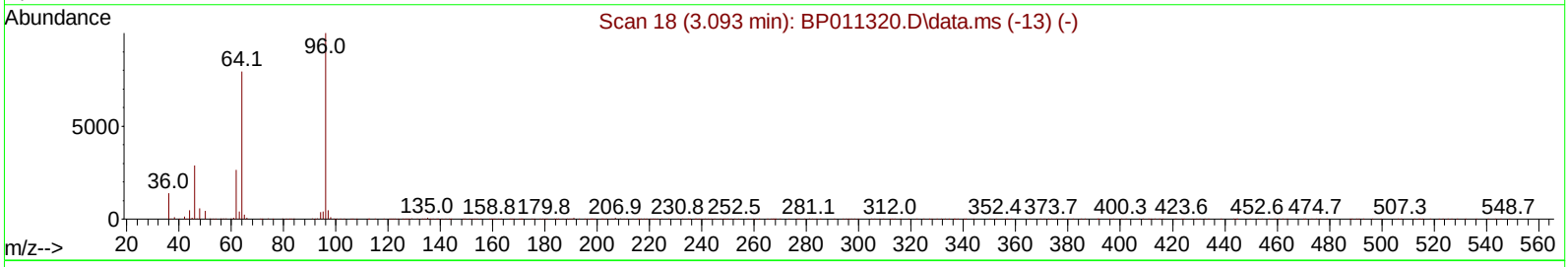

TIC: BP011325.D\data.ms

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

3.099min (+ 0.012) 4.04 ng/uL m

response 20847

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	61.00	74.54#
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	158226	20.000	ng/ul	0.00
20) Naphthalene-d8	10.381	136	674158	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.263	164	381383	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.033	188	769399	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	730911	20.000	ng/ul	0.00
88) Perylene-d12	23.468	264	732912	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.099	96	20847m	4.039	ng/uL	0.01
4) Pyridine-d5	3.511	84	92430	7.099	ng/ul	0.01
7) Phenol-d5	6.775	99	55090	3.934	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	197733	19.916	ng/ul	0.00
11) 2-Chlorophenol-d4	7.122	132	172643	15.714	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	96286	8.758	ng/ul	0.00
21) Nitrobenzene-d5	8.752	128	100272	21.855	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	92030	21.908	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.004	165	165853	19.551	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	119897	8.084	ng/ul	0.00
46) Dimethylphthalate-d6	13.687	166	673856	25.465	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	735396	23.086	ng/ul	0.00
54) 4-Nitrophenol-d4	14.492	143	11768	2.435	ng/ul	0.00
60) Fluorene-d10	15.263	176	545633	24.111	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.398	200	76357	21.366	ng/ul	0.00
73) Anthracene-d10	17.133	188	900397	26.036	ng/ul	0.00
81) Pyrene-d10	19.392	212	1094720	27.185	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.321	264	1033726	27.968	ng/ul	0.00

Target Compounds Qvalue

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

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