Quantitation Report (Qedit)

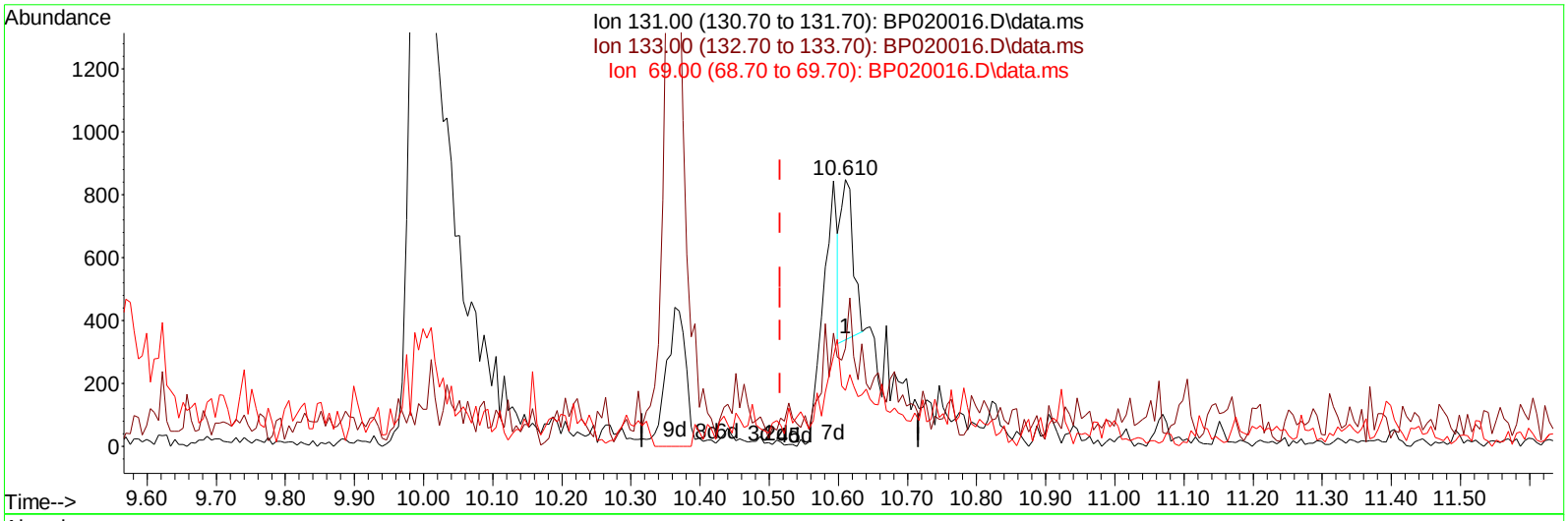
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032824\
Data File : BP020016.D
Acq On : 28 Mar 2024 18:47
Operator : MA/JU
Sample : PB159842BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

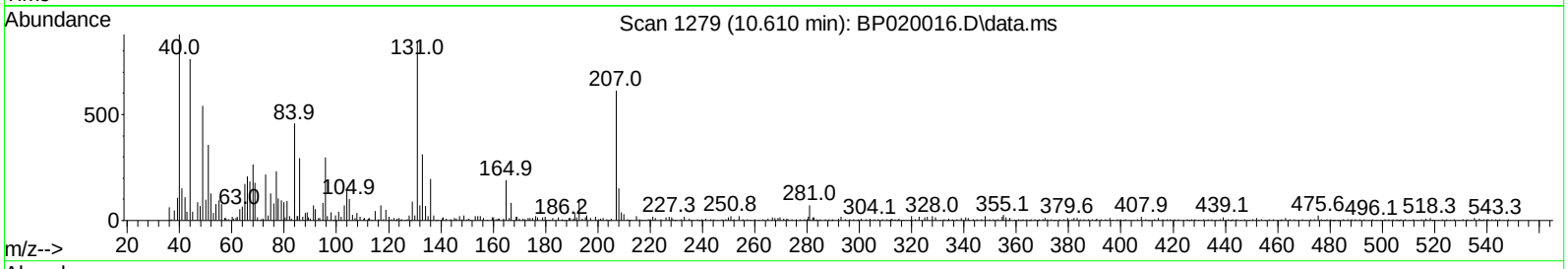
Quant Time: Mar 28 23:15:21 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 28 14:08:48 2024
Response via : Initial Calibration

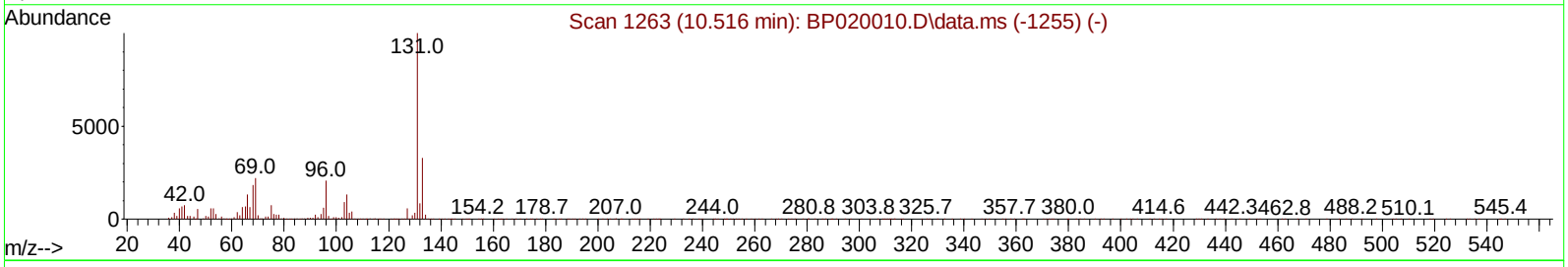
Instrument :
BNA_P
ClientSampleId :
SBLK842

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 03/29/2024
Supervised By :Jagrut Upadhyay 03/29/2024







TIC: BP020016.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.610min (+ 0.094) 0.06 ng/ul

response 625

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	36.95
69.00	22.40	21.02
0.00	0.00	0.00

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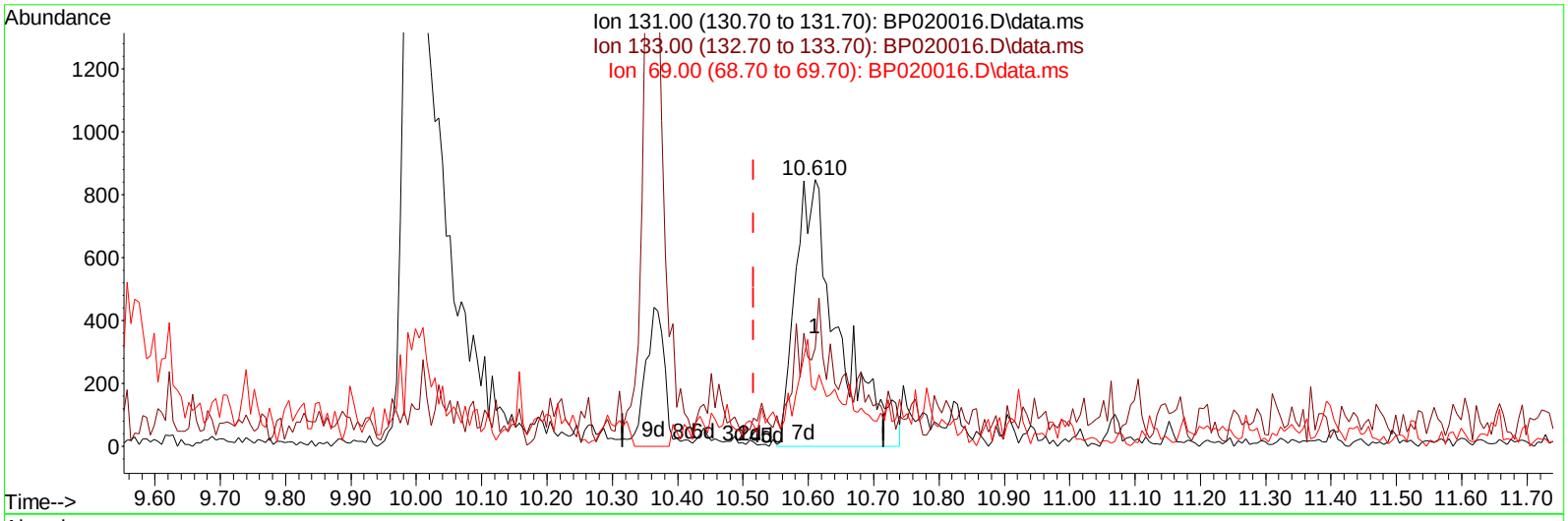

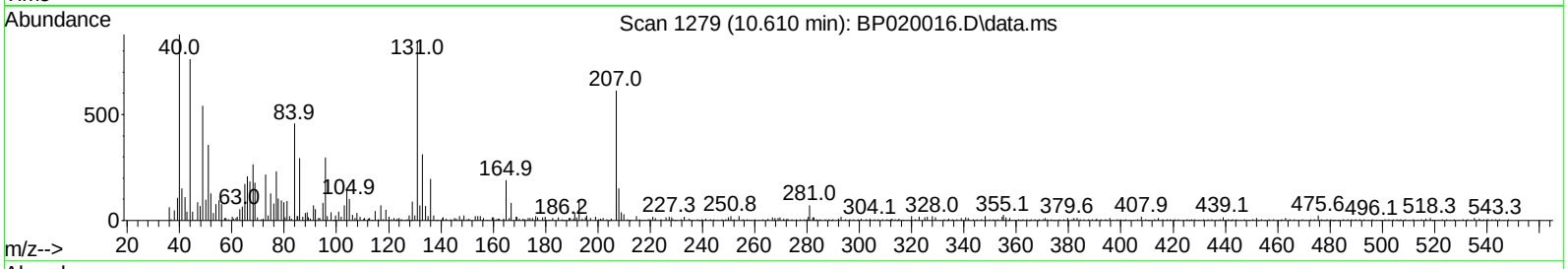

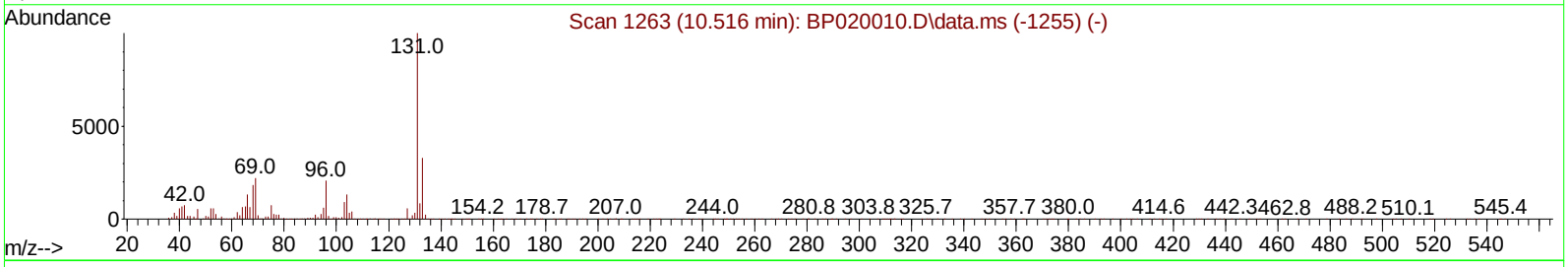

TIC: BP020016.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.610min (+ 0.094) 0.37 ng/ul m

response 3926

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	36.95
69.00	22.40	21.02
0.00	0.00	0.00

Quantitation Report (QT Reviewed)

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032824\
Data File : BP020016.D
Acq On : 28 Mar 2024 18:47
Operator : MA/JU
Sample : PB159842BL
Misc :
ALS Vial : 7 Sample Multiplier: 1

Quant Time: Mar 28 23:16:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Mar 28 14:08:48 2024
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Instrument :
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.617	152	126650	20.000	ng/ul	0.00
20) Naphthalene-d8	10.358	136	525597	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.251	164	355556	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.063	188	712505	20.000	ng/ul	-0.01
79) Chrysene-d12	21.539	240	543130	20.000	ng/ul	0.00
88) Perylene-d12	24.810	264	546107	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	19963	6.906	ng/uL	0.00
4) Pyridine-d5	3.646	84	101919	11.677	ng/ul	0.00
7) Phenol-d5	6.805	99	328830	30.785	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.964	67	231739	33.300	ng/ul	0.00
11) 2-Chlorophenol-d4	7.158	132	244395	31.859	ng/ul	0.00
15) 4-Methylphenol-d8	8.316	113	255339	30.898	ng/ul	0.01
21) Nitrobenzene-d5	8.758	128	124333	30.863	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	141450	31.382	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.999	165	237434	28.843	ng/ul	0.00
31) 4-Chloroaniline-d4	10.610	131	3926m	0.366	ng/ul	0.09
46) Dimethylphthalate-d6	13.663	166	812360	32.587	ng/ul	0.01
49) Acenaphthylene-d8	13.934	160	896181	30.775	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	101934	26.010	ng/ul	0.02
60) Fluorene-d10	15.269	176	695047	32.730	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.404	200	120849	28.335	ng/ul	-0.02
73) Anthracene-d10	17.181	188	1049242	33.267	ng/ul	0.01
81) Pyrene-d10	19.569	212	1144834	35.926	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.592	264	863259	32.904	ng/ul	-0.02
Target Compounds						Qvalue

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

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