Quantitation Report (Qedit)

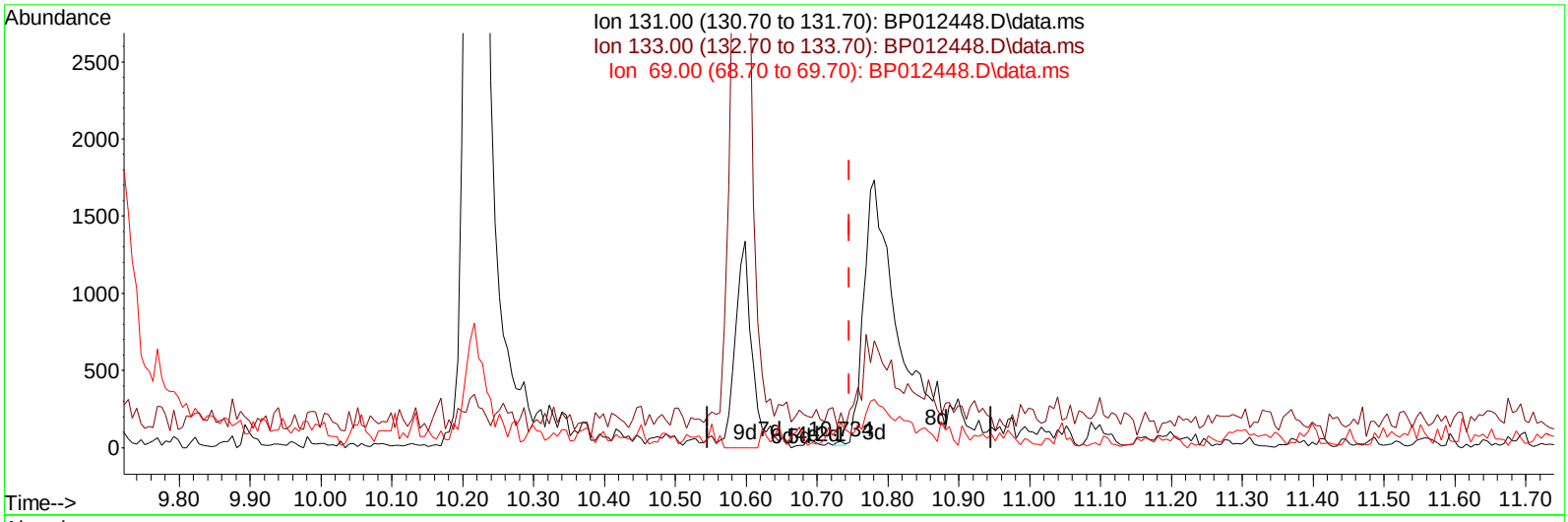
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110222\
Data File : BP012448.D
Acq On : 02 Nov 2022 11:53
Operator : CG/JU
Sample : PB148675BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

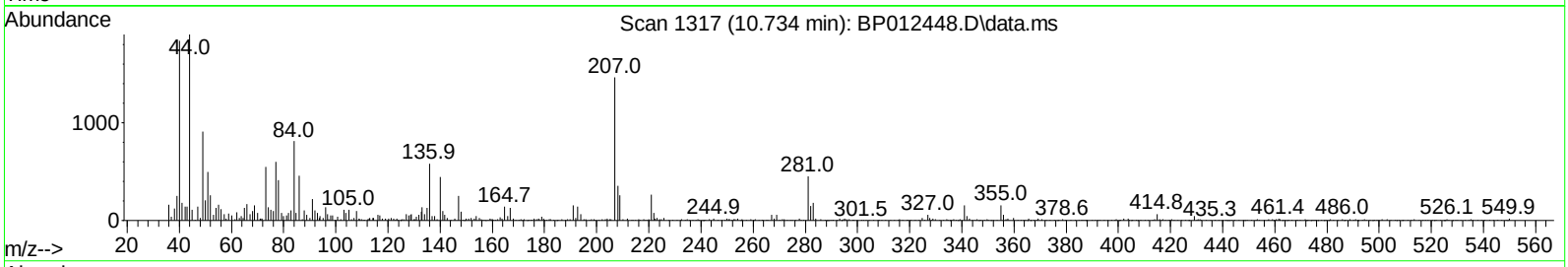
Quant Time: Nov 02 23:13:37 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 01 01:03:58 2022
Response via : Initial Calibration

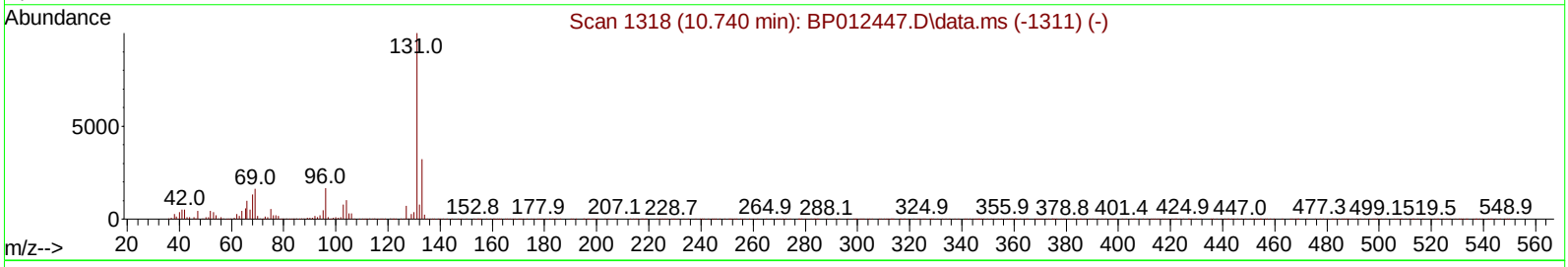
Instrument :
BNA_P
ClientSampleId :
SBLK675

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/03/2022
Supervised By :mohammad ahmed 11/03/2022







TIC: BP012448.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.734min (-0.012) 0.00 ng/ul

response 21

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	572.50#
69.00	16.50	395.00#
0.00	0.00	0.00

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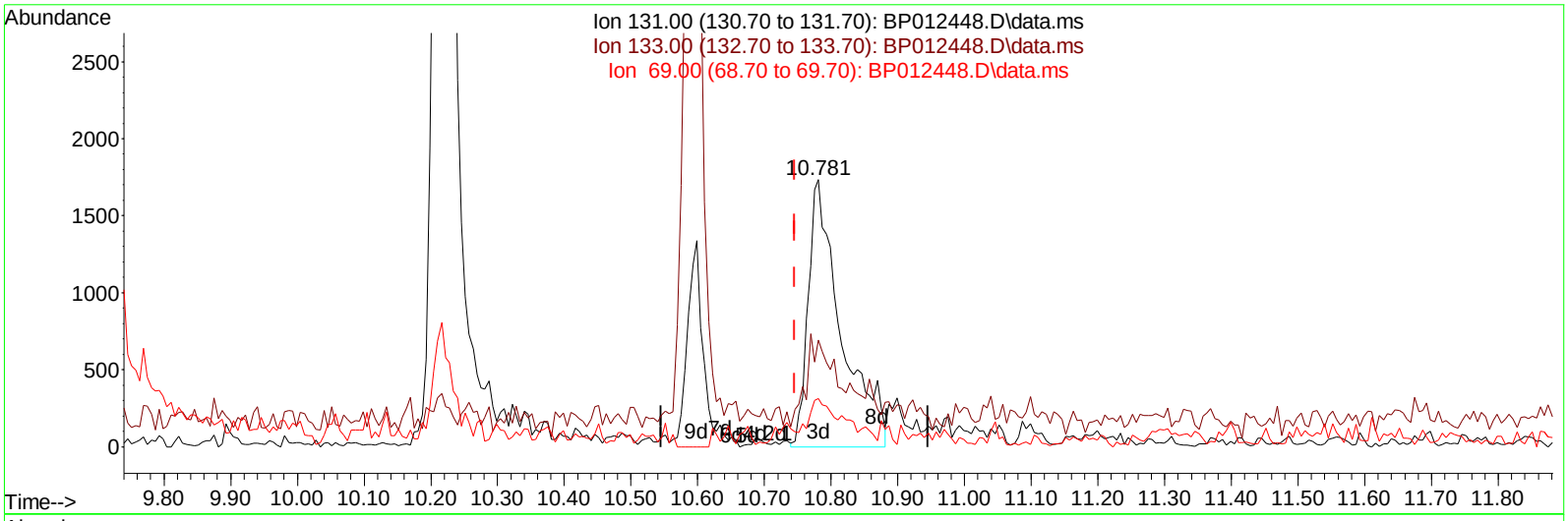

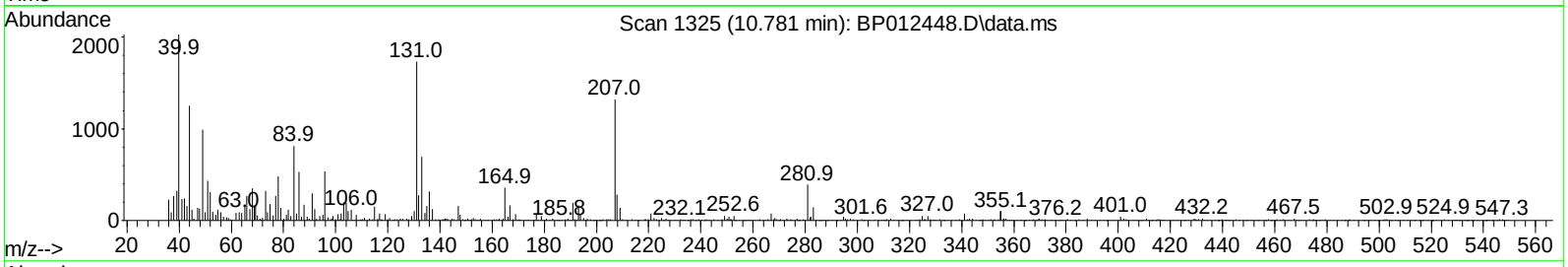

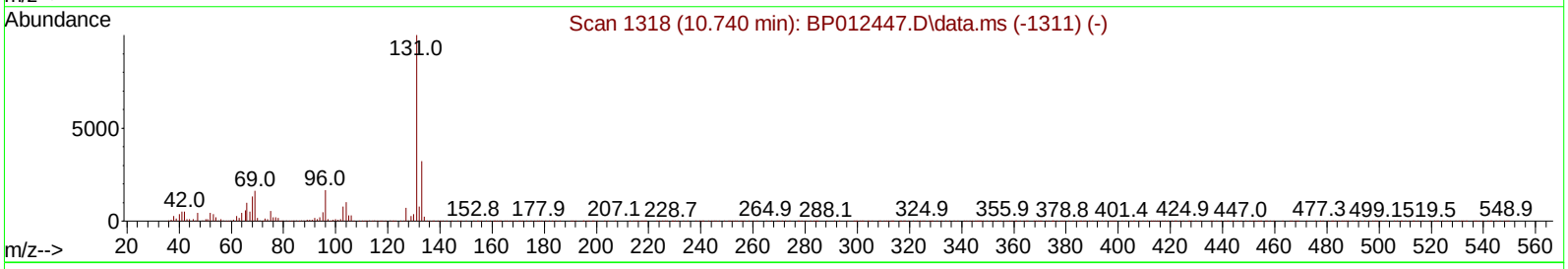

TIC: BP012448.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.781min (+ 0.035) 0.22 ng/ul m

response 5945

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	39.99#
69.00	16.50	9.06#
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	7.793	152	275246	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	1254380	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	861082	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1913465	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	1663864	20.000	ng/ul	0.00
88) Perylene-d12	23.686	264	1471392	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	41996	6.143	ng/uL	0.00
4) Pyridine-d5	3.670	84	74811	3.810	ng/ul	0.01
7) Phenol-d5	6.963	99	745829	30.211	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	451576	30.641	ng/ul	0.00
11) 2-Chlorophenol-d4	7.322	132	591883	32.831	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	608524	31.166	ng/ul	0.00
21) Nitrobenzene-d5	8.957	128	307378	37.787	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	355073	42.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.216	165	591499	32.327	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	5945m	0.218	ng/ul	0.04
46) Dimethylphthalate-d6	13.857	166	2067607	35.076	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	2254270	33.465	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	316918	28.407	ng/ul	0.00
60) Fluorene-d10	15.439	176	1791762	35.503	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	205549	20.633	ng/ul	0.00
73) Anthracene-d10	17.304	188	2876525	34.889	ng/ul	0.00
81) Pyrene-d10	19.545	212	3208049	38.403	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.533	264	2517669	35.019	ng/ul	-0.01
Target Compounds						Qvalue

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

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