

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019134.D
Acq On : 28 Dec 2023 20:31
Operator : MA/JU
Sample : 05920-11
Misc :
ALS Vial : 11 Sample Multiplier: 1

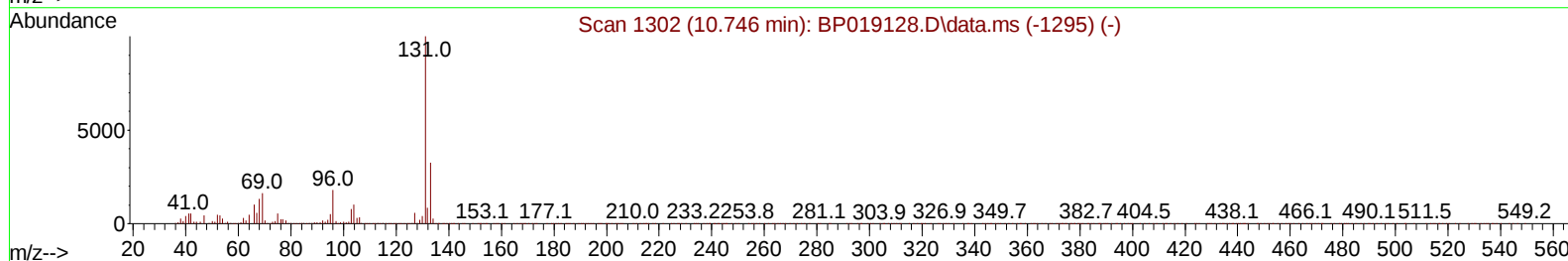
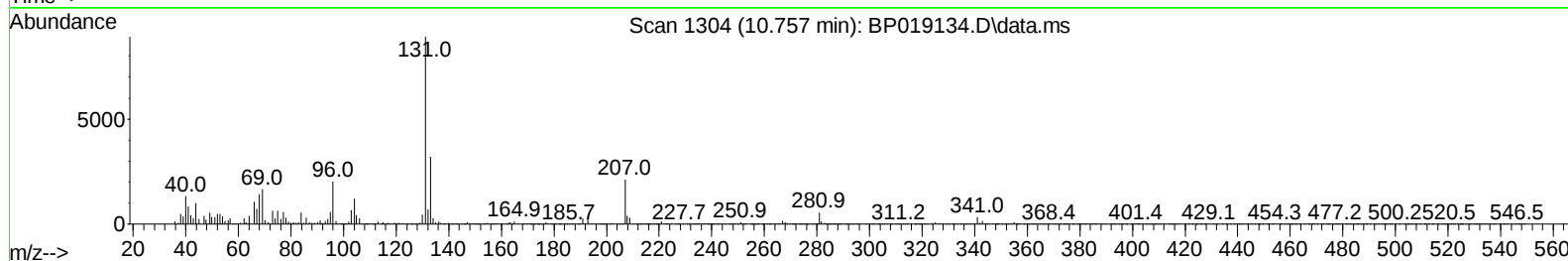
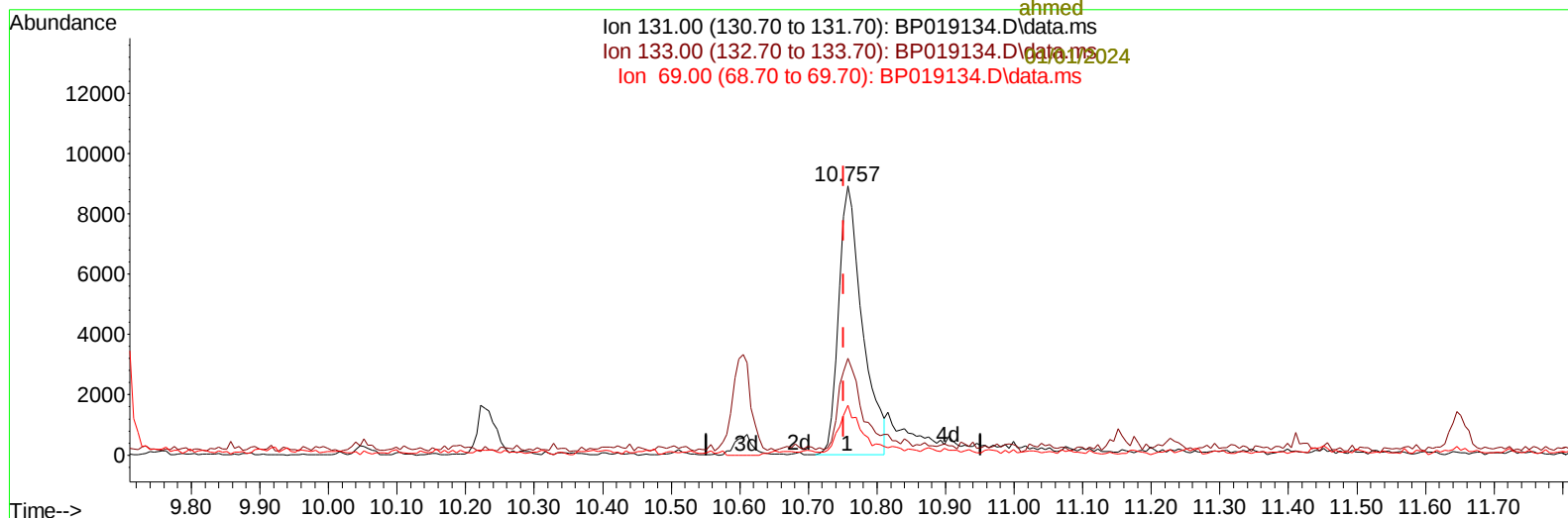
Instrument :
BNA_P
ClientSampleId :
GCF62

Manual IntegrationsAPPROVED

Quant Time: Dec 28 21:50:15 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 28 03:15:29 2023
Response via : Initial Calibration

Reviewed By :Yogesh
Patel

12/29/2023
Supervised By :mohammad
ahmed



TIC: BP019134.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.757min (+ 0.006) 1.26 ng/ul

response 21298

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.60	36.07
69.00	16.10	18.60
0.00	0.00	0.00

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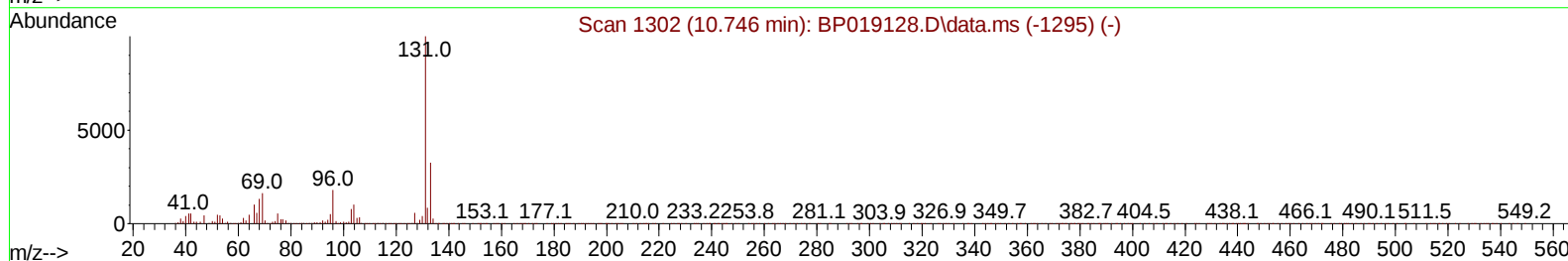
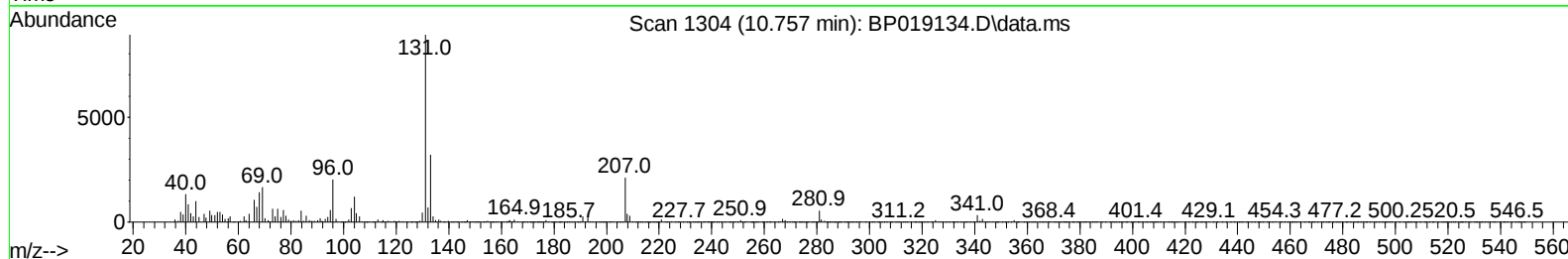
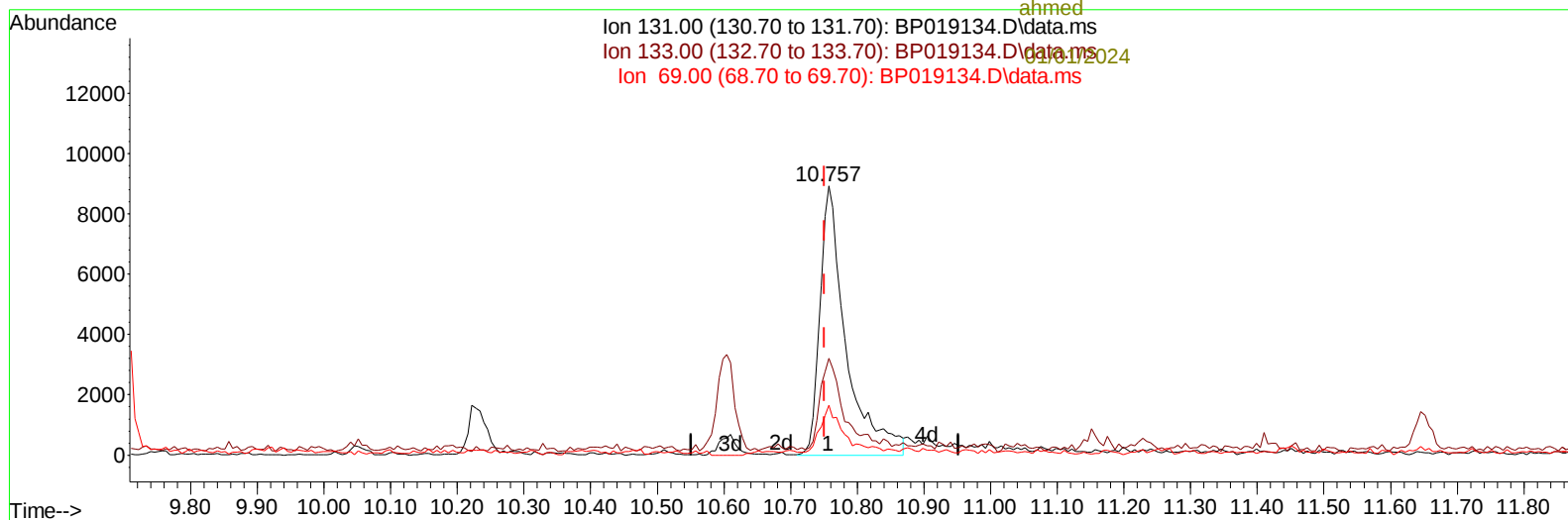
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TIC: BP019134.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.757min (+ 0.006) 1.44 ng/ul m

response 24244

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.60	36.07
69.00	16.10	18.60
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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Quant Time: Dec 28 21:58:54 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 28 03:15:29 2023
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----01/01/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.805	152	190936	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	726066	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	430506	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	978222	20.000	ng/ul	0.00
79) Chrysene-d12	21.310	240	1098003	20.000	ng/ul	0.00
88) Perylene-d12	23.674	264	1288194	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	12653	2.263	ng/uL	0.00
4) Pyridine-d5	3.664	84	32591	2.323	ng/ul	0.00
7) Phenol-d5	6.981	99	77333	4.516	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	130678	13.192	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	132257	9.581	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	58510	4.389	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	74716	11.904	ng/ul	0.00
24) 2-Nitrophenol-d4	9.693	143	78170	11.524	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	91963	6.714	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	24244m	1.439	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	470235	12.366	ng/ul	0.00
49) Acenaphthylene-d8	14.140	160	271462	6.579	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	25213	3.927	ng/ul	0.00
60) Fluorene-d10	15.445	176	422871	12.872	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	50709	7.907	ng/ul	-0.01
73) Anthracene-d10	17.304	188	372971	7.301	ng/ul	0.00
81) Pyrene-d10	19.551	212	99968	1.494	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.515	264	47176	0.637	ng/ul	-0.02
Target Compounds						
					Qvalue	
8) Phenol	7.010	94	27159	1.606	ng/ul	97
16) Acetophenone	8.634	105	134894	6.756	ng/ul	98
71) Pentachlorophenol	16.857	266	27439	3.145	ng/ul	96

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

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