

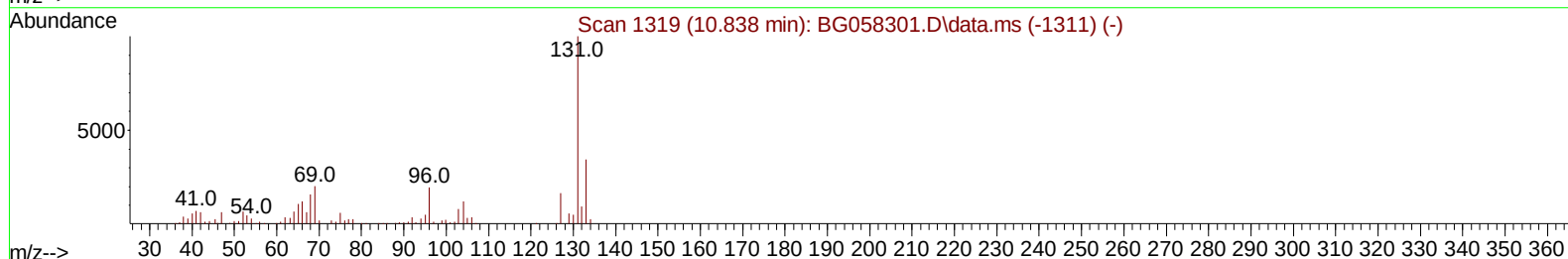
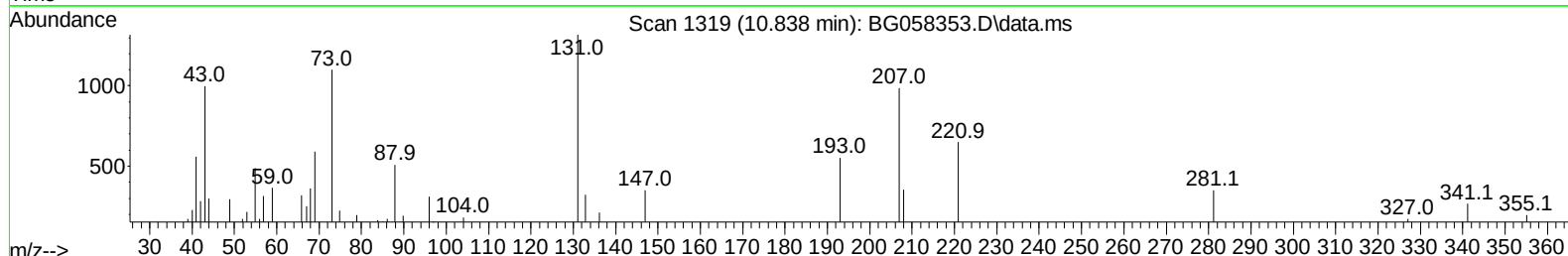
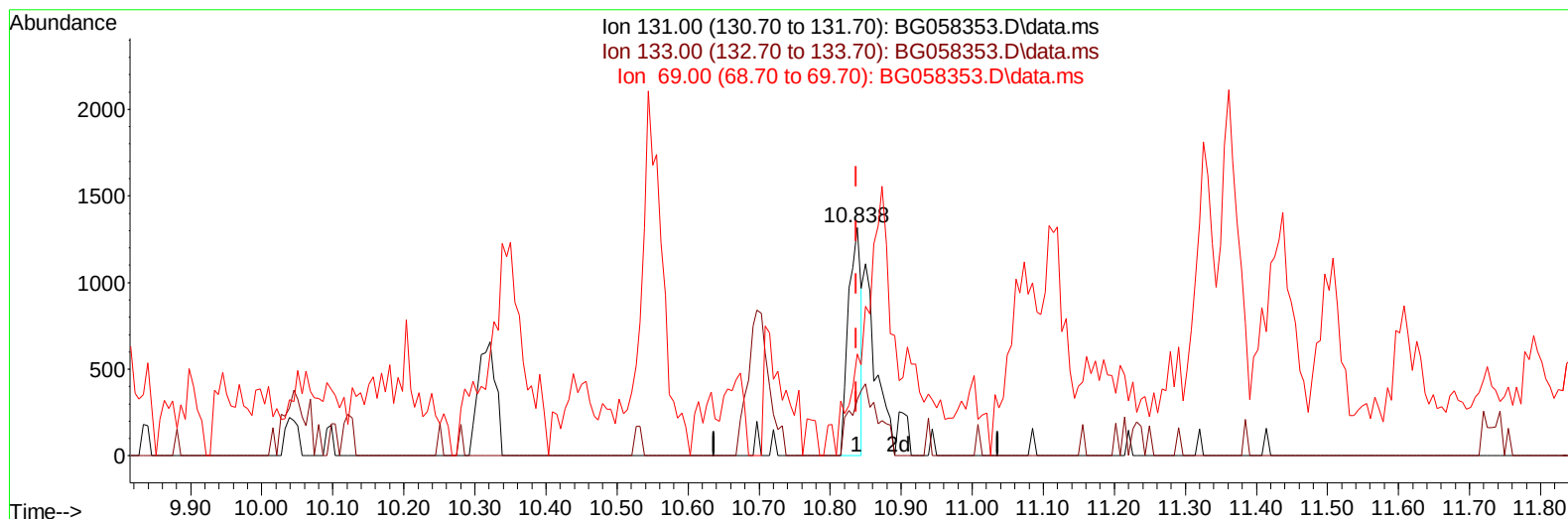
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071823\
Data File : BG058353.D
Acq On : 20 Jul 2023 9:18
Operator : CG/JU
Sample : 03452-17
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
A46P9

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/20/2023
Supervised By :Sohil Jodhani 07/20/2023

Quant Time: Jul 20 09:55:37 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jul 18 02:19:39 2023
Response via : Initial Calibration



TIC: BG058353.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.838min (+ 0.001) 0.39 ng/ul

response 1670

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.90	24.51#
69.00	18.60	44.76#
0.00	0.00	0.00

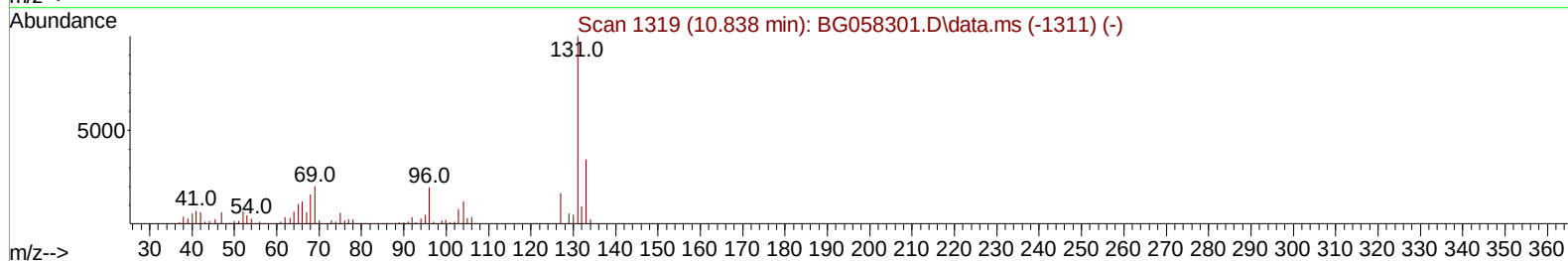
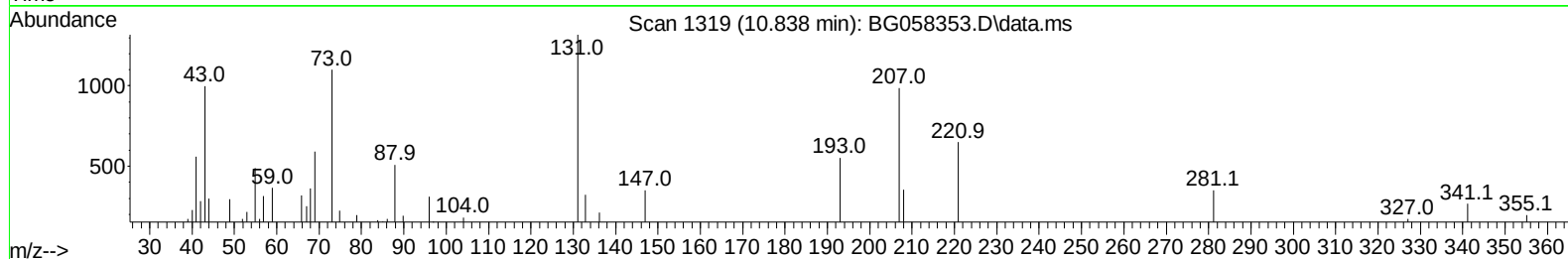
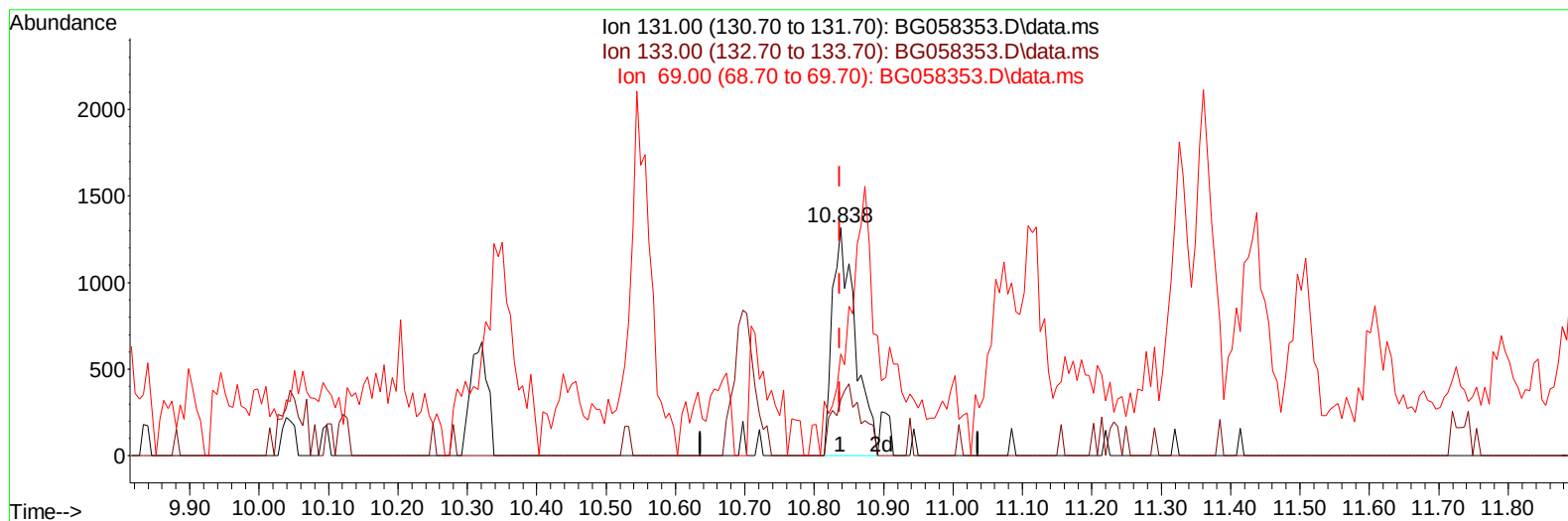
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(31) 4-Chloroaniline-d4 (S)

10.838min (+ 0.001) 0.70 ng/ul m

response 3019

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.90	24.51#
69.00	18.60	44.76#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 07/20/2023
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Quant Time: Jul 20 09:56:58 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG071723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jul 18 02:19:39 2023
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.894	152		39535	20.000	ng/ul	0.00
20) Naphthalene-d8	10.697	136		190248	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.534	164		137377	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.278	188		311709	20.000	ng/ul	0.00
79) Chrysene-d12	21.525	240		256748	20.000	ng/ul	0.00
88) Perylene-d12	24.657	264		289807	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.335	96		4215	3.924	ng/uL	0.00
4) Pyridine-d5	3.746	84		67640	21.216	ng/ul	0.00
7) Phenol-d5	7.060	99		60454	16.119	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.219	67		49183	20.982	ng/ul	0.00
11) 2-Chlorophenol-d4	7.430	132		46806	17.829	ng/ul	0.00
15) 4-Methylphenol-d8	8.611	113		1881	0.658	ng/ul	0.00
21) Nitrobenzene-d5	9.058	128		30128	21.127	ng/ul	0.00
24) 2-Nitrophenol-d4	9.780	143		33555	21.554	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.321	165		37048	12.414	ng/ul	0.00
31) 4-Chloroaniline-d4	10.838	131		3019m	0.698	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166		207858	19.621	ng/ul	0.00
49) Acenaphthylene-d8	14.228	160		215313	19.395	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143		27190	16.995	ng/ul	0.00
60) Fluorene-d10	15.527	176		190530	21.724	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.644	200		24277	15.287	ng/ul	0.00
73) Anthracene-d10	17.372	188		123010	9.133	ng/ul	-0.01
81) Pyrene-d10	19.669	212		324139	21.375	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.440	264		100099	7.146	ng/ul	-0.01
Target Compounds							
						Qvalue	
16) Acetophenone	8.717	105		16348	3.356	ng/ul	86
80) Fluoranthene	19.328	202		19499	1.122	ng/ul	98
86) Bis(2-ethylhexyl)phtha...	21.408	149		20306	2.060	ng/ul	98

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

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