

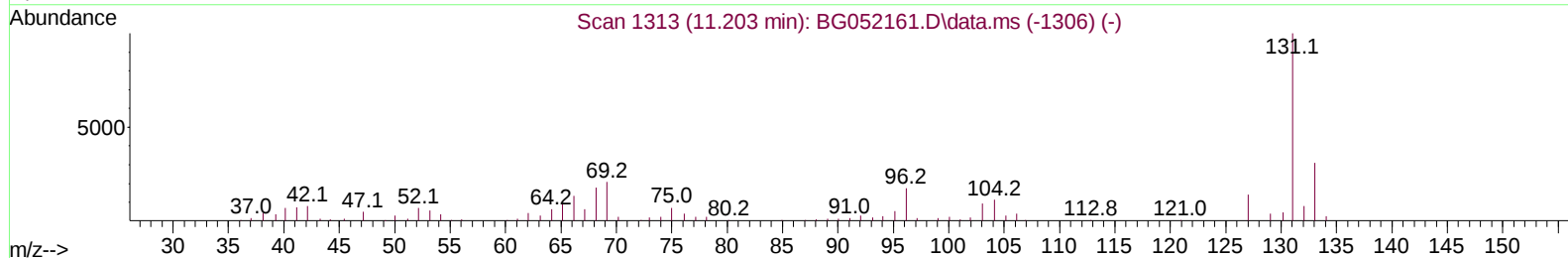
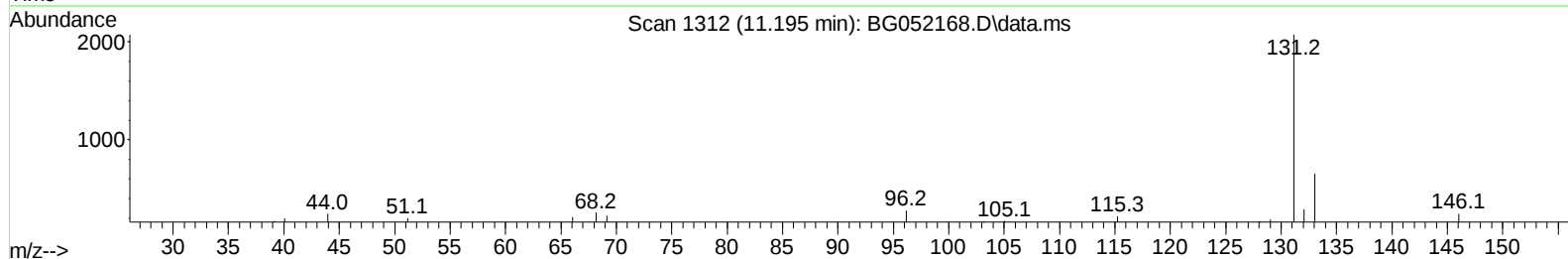
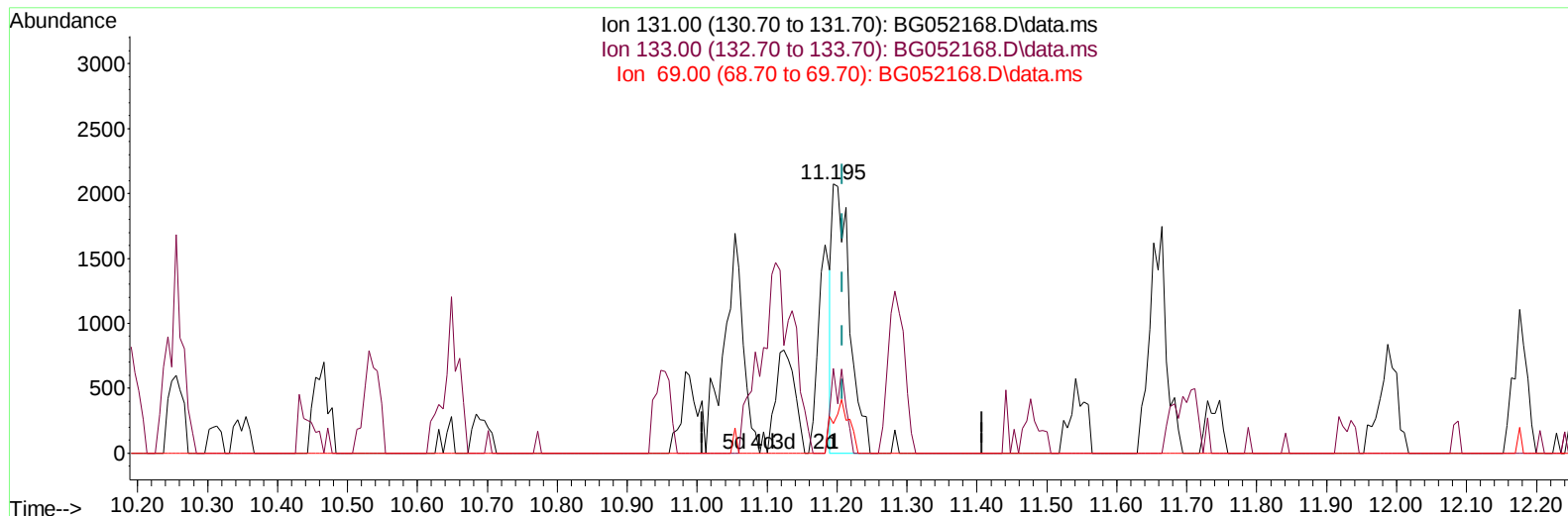
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052168.D
 Acq On : 27 Jan 2022 3:34
 Operator : CG/JU
 Sample : N1230-03DL 10X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0Q24DL

Manual IntegrationsAPPROVED

Quant Time: Jan 27 04:16:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/27/2022
 Supervised By :Yogesh Patel 01/28/2022



TIC: BG052168.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.195min (-0.013) 1.32 ng/u1

response 3598

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.20	31.56
69.00	21.30	10.96#
0.00	0.00	0.00

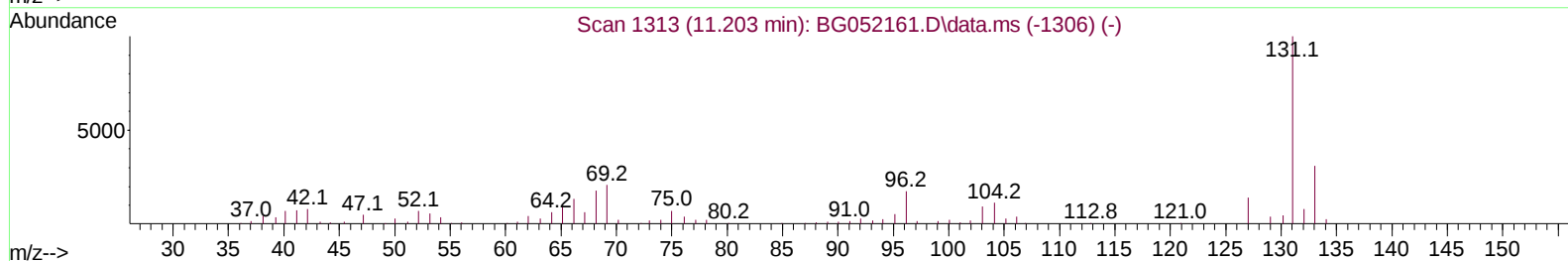
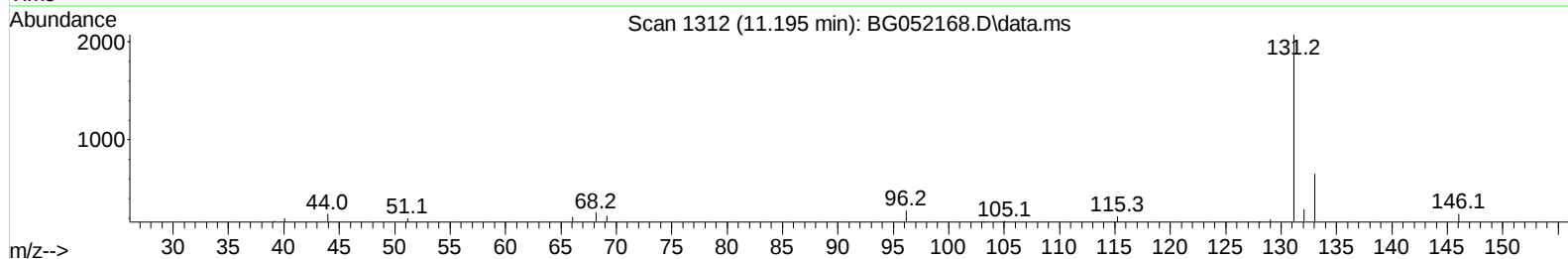
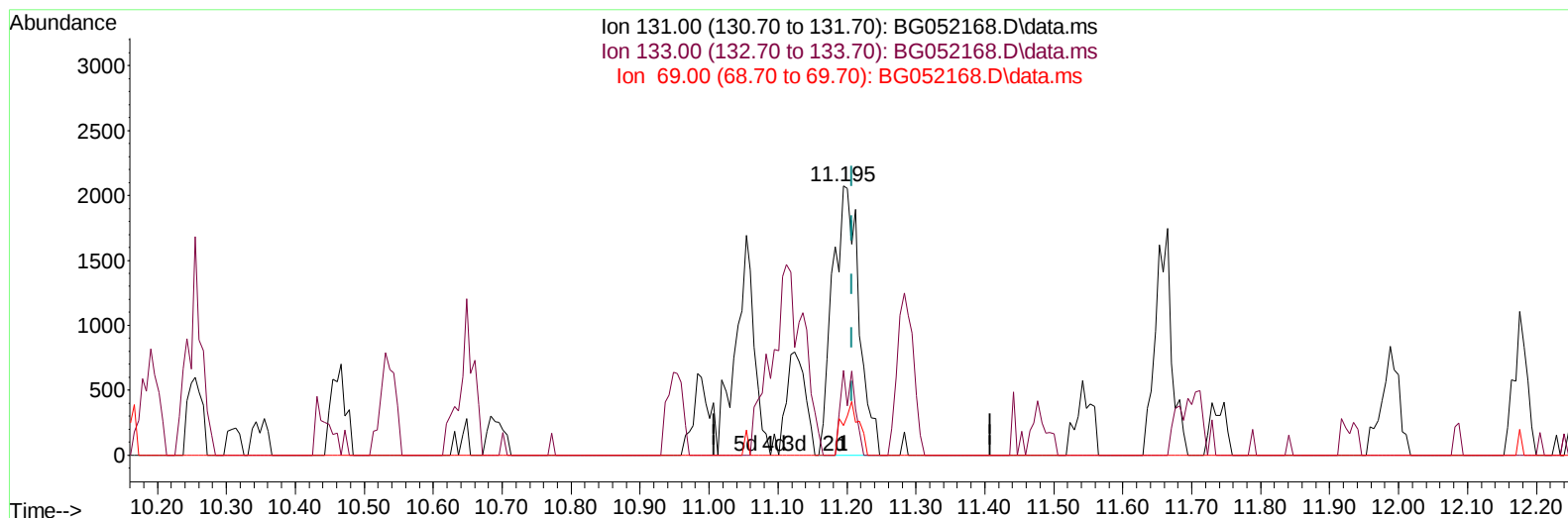
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TIC: BG052168.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.195min (-0.013) 2.02 ng/ul m

response 5499

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.20	31.56
69.00	21.30	10.96#
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	8.240	152	26174	20.000	ng/ul	0.00
20) Naphthalene-d8	11.077	136	114258	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.884	164	86767	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.633	188	214126	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.945	240	227971	20.000	ng/ul	# 0.00
88) Perylene-d12	25.423	264	227176	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0d	0.000	ng/uL	
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.376	99	1623	0.634	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.541	67	4845	2.979	ng/ul	0.00
11) 2-Chlorophenol-d4	7.764	132	4544	2.679	ng/ul	0.00
15) 4-Methylphenol-d8	8.939	113	3110	1.562	ng/ul	0.00
21) Nitrobenzene-d5	9.409	128	2976	3.296	ng/ul	0.00
24) 2-Nitrophenol-d4	10.137	143	3045	3.050	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.690	165	5760	2.951	ng/ul	0.00
31) 4-Chloroaniline-d4	11.195	131	5499m	2.023	ng/ul	-0.01
46) Dimethylphthalate-d6	14.267	166	26998	3.764	ng/ul	0.00
49) Acenaphthylene-d8	14.573	160	31545	3.641	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.865	176	24112	3.856	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.977	200	2780	2.058	ng/ul	0.00
73) Anthracene-d10	17.727	188	41830	4.118	ng/ul	0.00
81) Pyrene-d10	20.006	212	51888	3.955	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.176	264	47463	3.969	ng/ul	0.00
Target Compounds						
					Qvalue	
52) Acenaphthene	14.948	153	7498	1.322	ng/ul	90
58) 2,3,4,6-Tetrachlorophenol	15.507	232	5750	3.032	ng/ul#	86
71) Pentachlorophenol	17.281	266	109221	70.347	ng/ul	94

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

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