

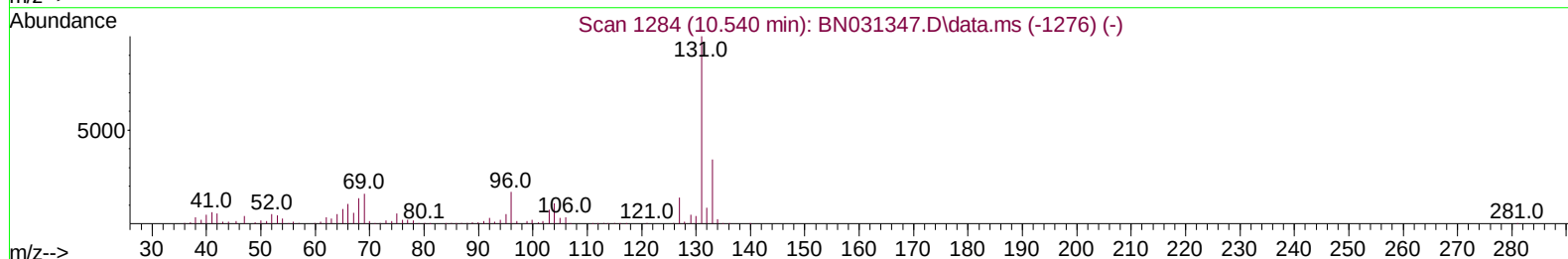
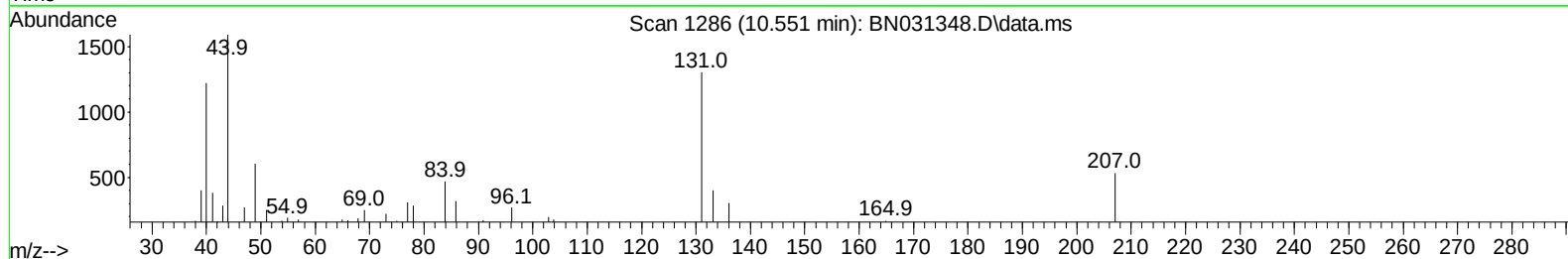
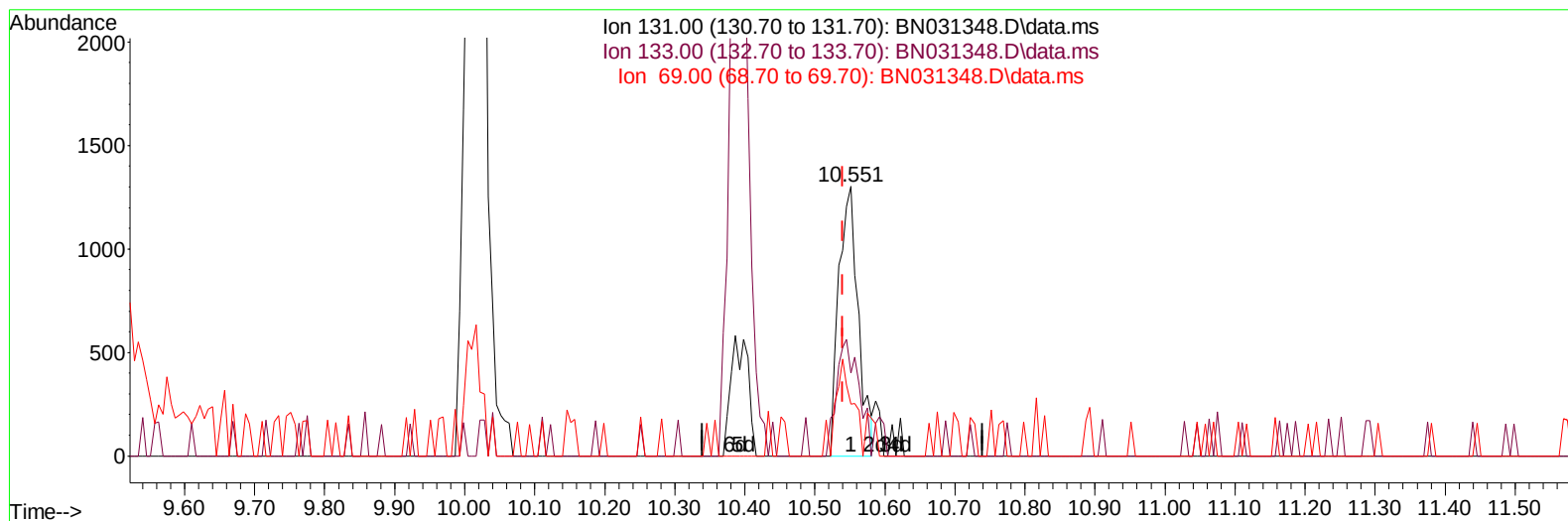
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050124\
Data File : BN031348.D
Acq On : 01 May 2024 13:38
Operator : MA/JU
Sample : PB160361BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK361

Manual IntegrationsAPPROVED

Quant Time: May 01 14:07:18 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed May 01 05:16:42 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/01/2024
Supervised By :mohammad ahmed 05/02/2024



TIC: BN031348.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.551min (+ 0.012) 0.18 ng/ul

response 2531

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.30	30.88
69.00	15.50	19.28#
0.00	0.00	0.00

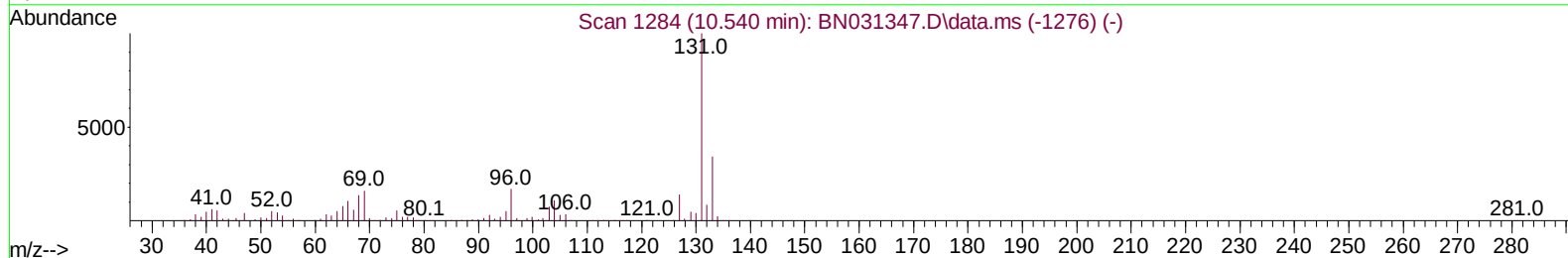
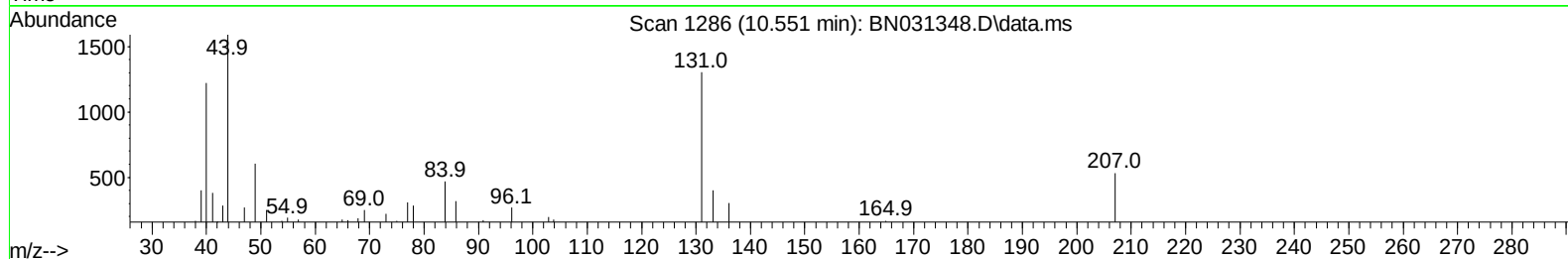
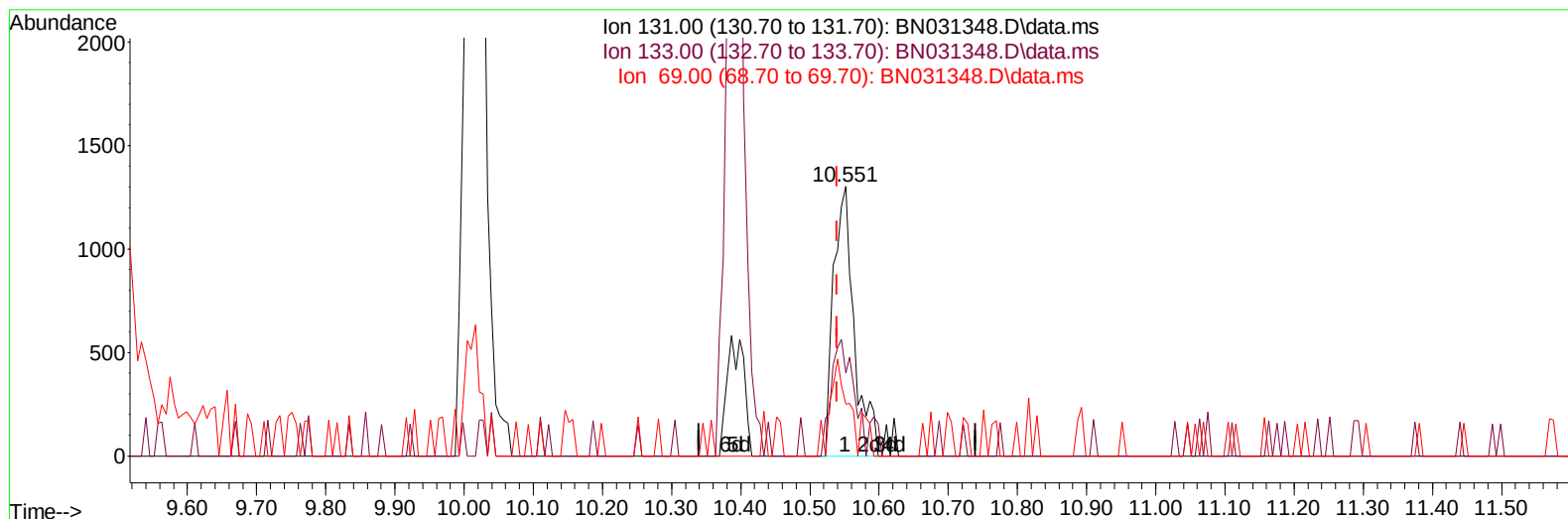
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Misc :
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Instrument :
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Manual IntegrationsAPPROVED

Quant Time: May 01 14:07:18 2024
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TIC: BN031348.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.551min (+ 0.012) 0.19 ng/ul m

response 2702

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.30	30.88
69.00	15.50	19.28#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN050124\
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 Operator : MA/JU
 Sample : PB160361BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK361

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 05/01/2024
 Supervised By :mohammad ahmed 05/02/2024

Quant Time: May 01 14:08:45 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN042924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 01 05:16:42 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.610	152	173610	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136	673899	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	408420	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.010	188	833546	20.000	ng/ul	0.00
79) Chrysene-d12	21.251	240	804619	20.000	ng/ul	0.00
88) Perylene-d12	24.145	264	926838	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.134	96	23266	6.239	ng/uL	0.00
4) Pyridine-d5	3.546	84	165593	15.507	ng/ul	0.01
7) Phenol-d5	6.787	99	333414	23.866	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.957	67	222945	27.256	ng/ul	0.00
11) 2-Chlorophenol-d4	7.146	132	310266	27.761	ng/ul	0.00
15) 4-Methylphenol-d8	8.316	113	234671	20.270	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128	148416	30.559	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	160130	30.635	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	269228	27.737	ng/ul	0.00
31) 4-Chloroaniline-d4	10.551	131	2702m	0.192	ng/ul	0.01
46) Dimethylphthalate-d6	13.675	166	852482	31.660	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	877268	28.080	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143	160262	32.797	ng/ul	0.00
60) Fluorene-d10	15.257	176	709404	30.853	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.386	200	124540	29.054	ng/ul	0.00
73) Anthracene-d10	17.110	188	1113321	32.363	ng/ul	0.00
81) Pyrene-d10	19.416	212	1337794	27.300	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.951	264	1173160	27.341	ng/ul	0.02

Target Compounds Qvalue

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

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