

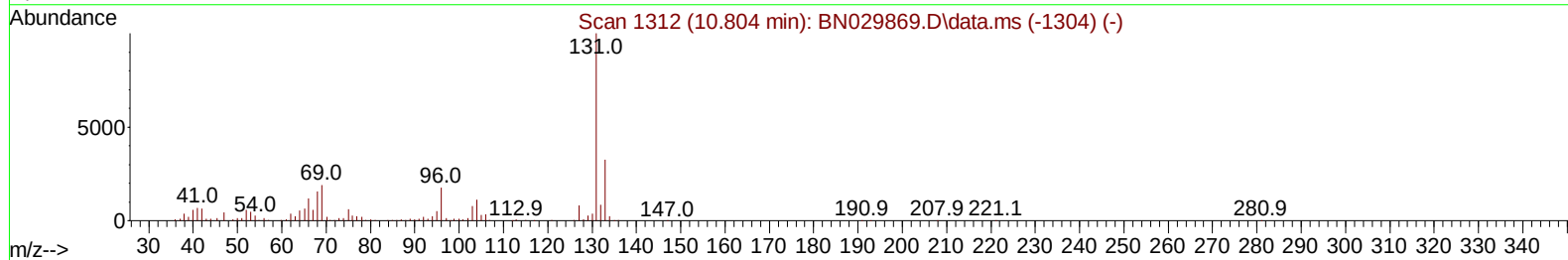
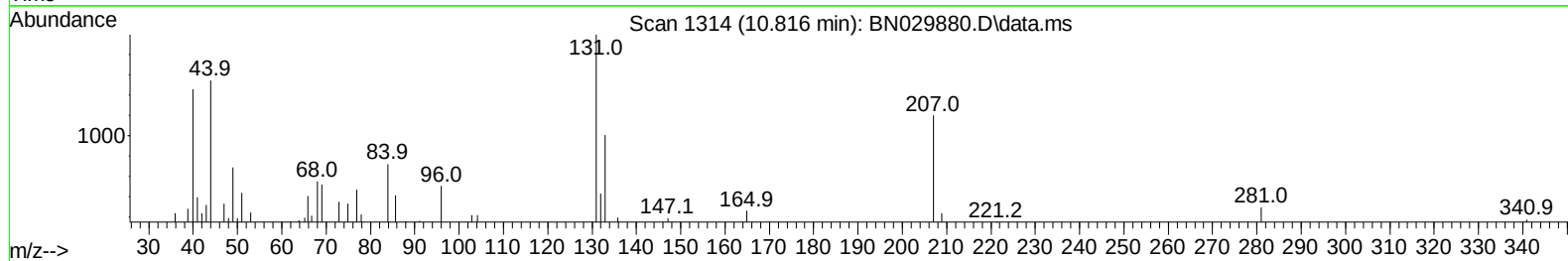
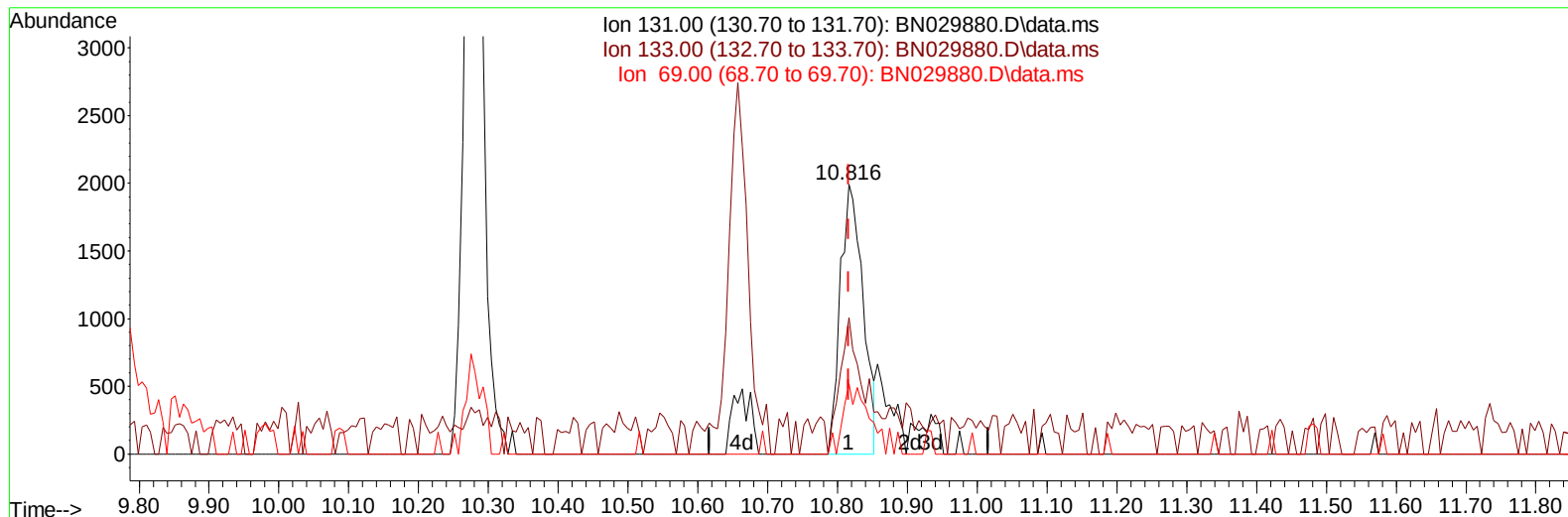
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021924\
Data File : BN029880.D
Acq On : 20 Feb 2024 17:28
Operator : MA/JU
Sample : P1455-13
Misc :
ALS Vial : 50 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
C0AF2

Manual IntegrationsAPPROVED

Quant Time: Feb 20 22:49:41 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Feb 13 22:50:45 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 02/21/2024
Supervised By :mohammad ahmed 02/21/2024



TIC: BN029880.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.816min (-0.000) 0.36 ng/u1

response 4486

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.10	50.48#
69.00	19.70	26.17#
0.00	0.00	0.00

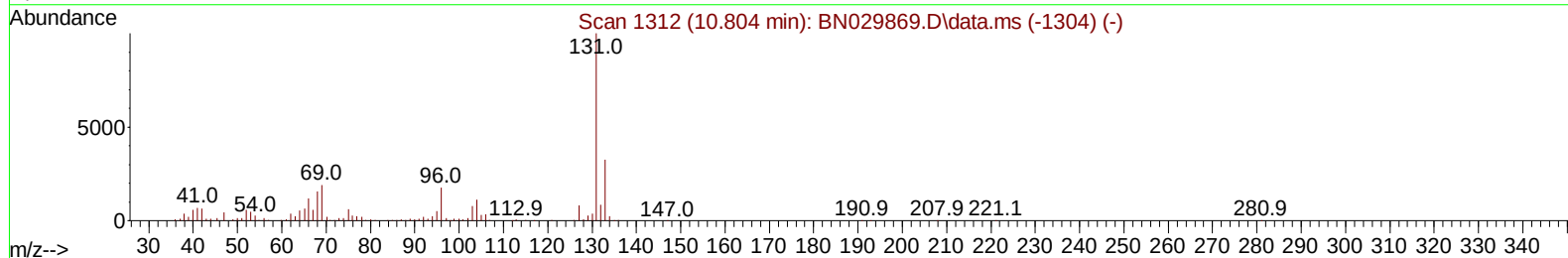
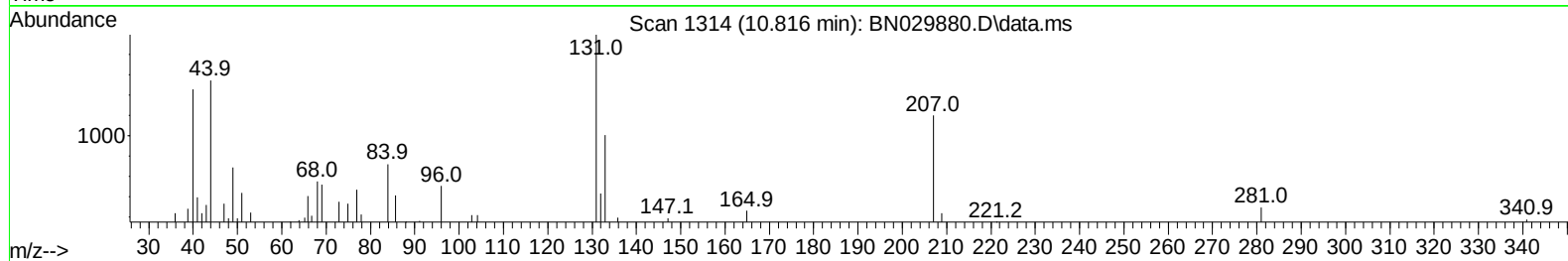
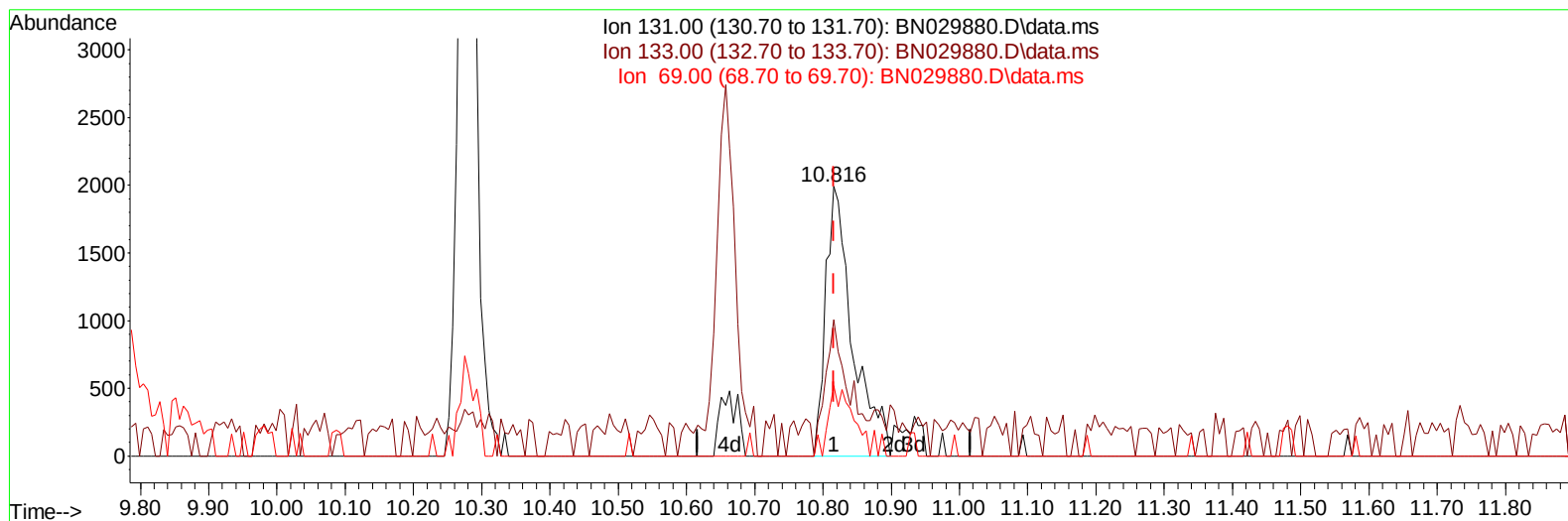
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TIC: BN029880.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.816min (-0.000) 0.43 ng/ul m

response 5453

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.10	50.48#
69.00	19.70	26.17#
0.00	0.00	0.00

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Quant Time: Feb 20 22:51:21 2024
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 QLast Update : Tue Feb 13 22:50:45 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.863	152	133163	20.000	ng/ul	0.00
20) Naphthalene-d8	10.657	136	551981	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.504	164	298602	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.257	188	615053	20.000	ng/ul	-0.01
79) Chrysene-d12	21.457	240	520320	20.000	ng/ul	-0.01
88) Perylene-d12	23.816	264	598648	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.317	96	17147	4.235	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.028	99	167162	13.421	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.205	67	321649	38.694	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.387	132	359267	39.804	ng/ul	-0.01
15) 4-Methylphenol-d8	8.569	113	263684	28.386	ng/ul	-0.02
21) Nitrobenzene-d5	9.028	128	185841	43.379	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.746	143	177937	47.690	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.281	165	317906	43.167	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.816	131	5453m	0.434	ng/ul	0.00
46) Dimethylphthalate-d6	13.928	166	979918	44.079	ng/ul	0.00
49) Acenaphthylene-d8	14.198	160	1112650	42.144	ng/ul	0.00
54) 4-Nitrophenol-d4	14.716	143	33448	7.927	ng/ul	0.00
60) Fluorene-d10	15.498	176	833389	44.066	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.628	200	104916	35.008	ng/ul	0.00
73) Anthracene-d10	17.357	188	1299437	45.080	ng/ul	0.00
81) Pyrene-d10	19.657	212	1436511	46.005	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.674	264	1358606	45.354	ng/ul	0.00

Target Compounds Qvalue

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

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