

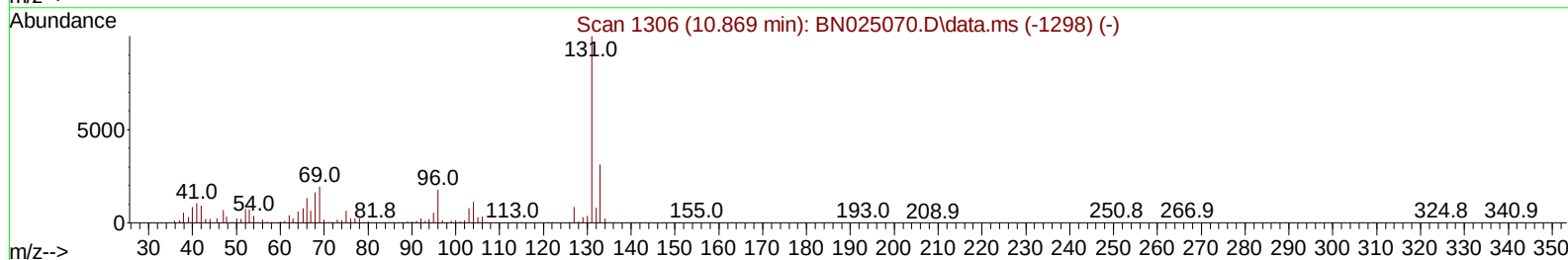
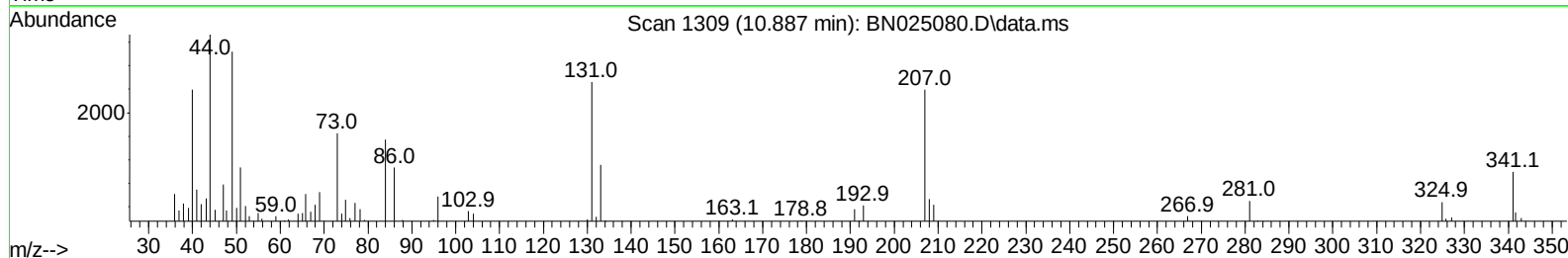
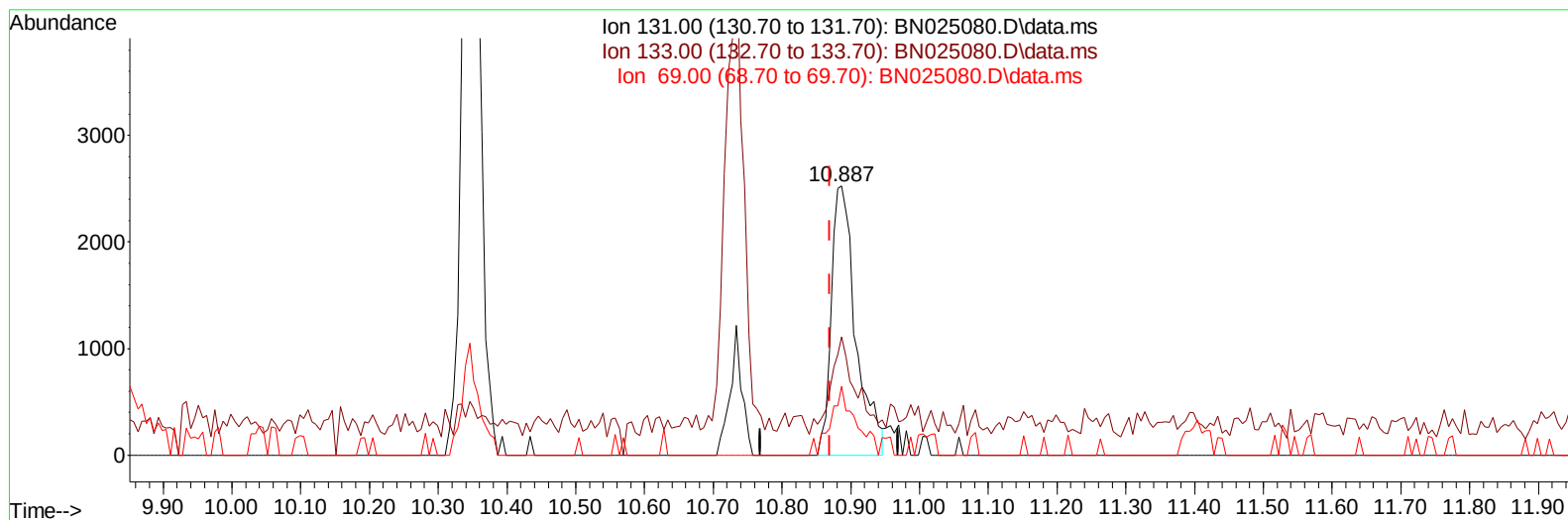
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN042223\
Data File : BN025080.D
Acq On : 22 Apr 2023 17:31
Operator : CG/JU
Sample : 02378-16
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
COMK9

Manual IntegrationsAPPROVED

Quant Time: Apr 23 00:41:17 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041023.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Apr 23 00:37:14 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/24/2023
Supervised By :Jagrut Upadhyay 04/24/2023



TIC: BN025080.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.887min (+ 0.018) 0.27 ng/ul

response 6255

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.00	43.92#
69.00	19.40	25.64#
0.00	0.00	0.00

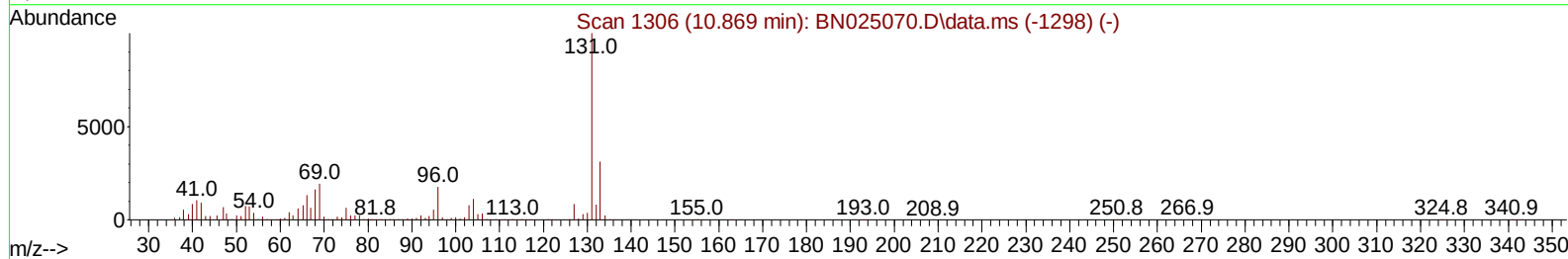
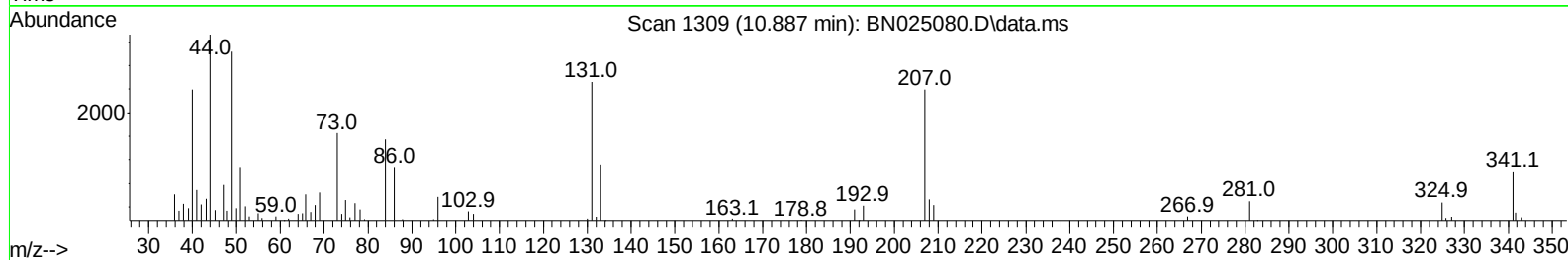
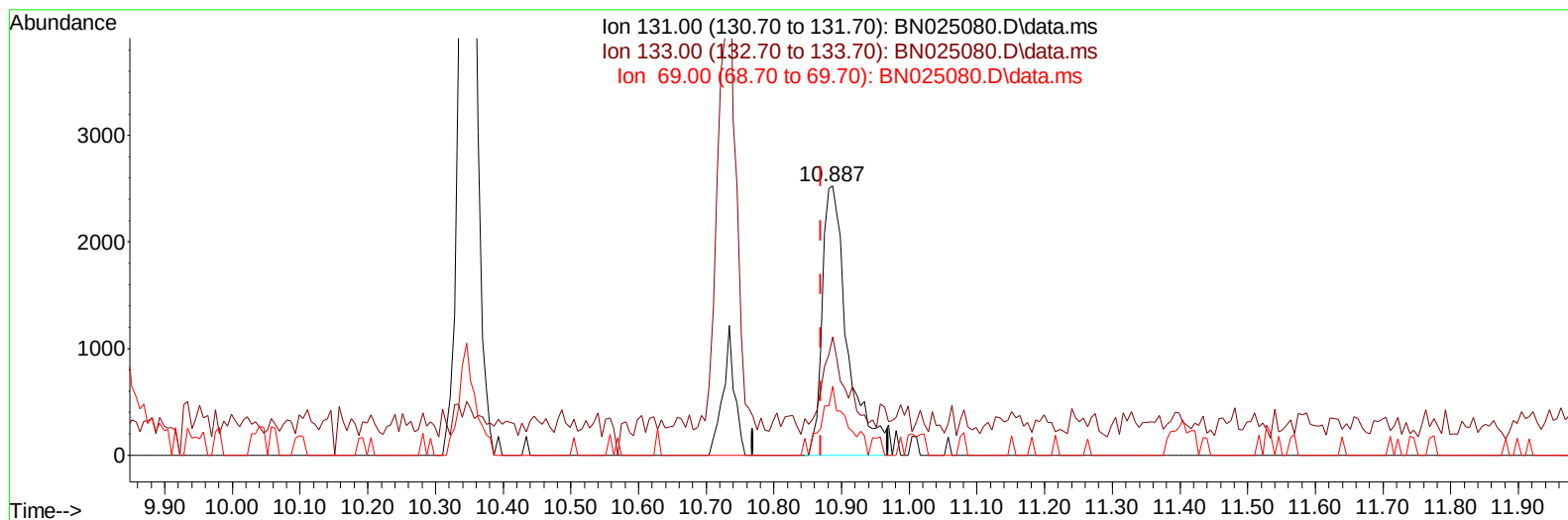
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TIC: BN025080.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.887min (+ 0.018) 0.29 ng/ul m

response 6617

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.00	43.92#
69.00	19.40	25.64#
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	7.916	152	210373	20.000	ng/ul	0.00
20) Naphthalene-d8	10.728	136	965714	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.563	164	635246	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.310	188	1440339	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	1485959	20.000	ng/ul	0.00
88) Perylene-d12	23.945	264	1746569	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.293	96	24655	4.460	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.075	99	221607	10.805	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.246	67	400869	29.222	ng/ul	0.00
11) 2-Chlorophenol-d4	7.446	132	408919	27.130	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	339980	20.880	ng/ul	0.00
21) Nitrobenzene-d5	9.087	128	216652	27.708	ng/ul	0.00
24) 2-Nitrophenol-d4	9.810	143	239682	28.300	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.346	165	414708	27.395	ng/ul	0.00
31) 4-Chloroaniline-d4	10.887	131	6617m	0.288	ng/ul	0.02
46) Dimethylphthalate-d6	13.975	166	1772832	36.000	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1696808	30.068	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	118088	11.698	ng/ul	0.00
60) Fluorene-d10	15.557	176	1470080	35.357	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	288669	30.796	ng/ul	0.00
73) Anthracene-d10	17.410	188	2557817	37.811	ng/ul	0.00
81) Pyrene-d10	19.704	212	3197071	37.052	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.792	264	3555477	38.933	ng/ul	0.01

Target Compounds Qvalue

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

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