

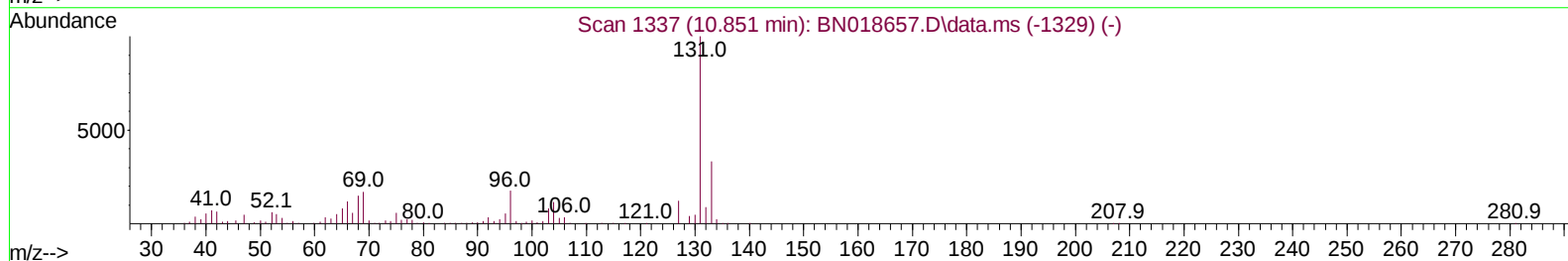
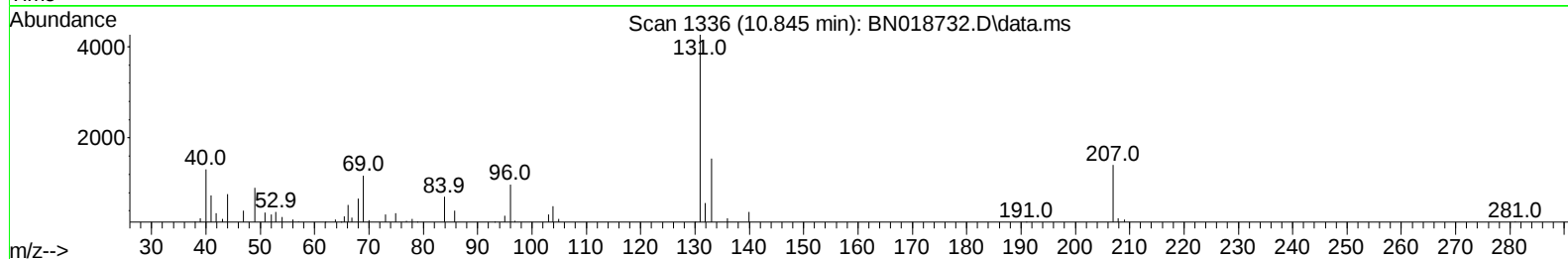
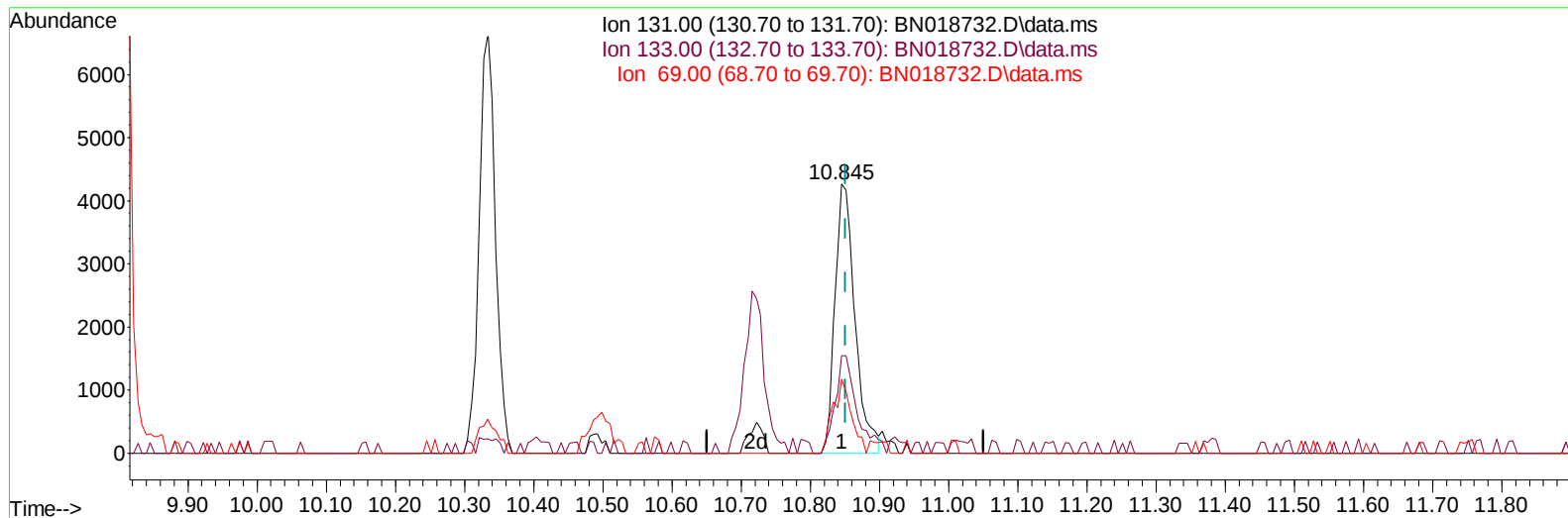
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022522\
Data File : BN018732.D
Acq On : 25 Feb 2022 15:54
Operator : CG/JU
Sample : N1656-08
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGQ15

Manual IntegrationsAPPROVED

Quant Time: Feb 25 18:01:01 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Feb 16 16:48:14 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/28/2022
Supervised By :mohammad ahmed 02/28/2022



TIC: BN018732.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.845min (-0.006) 0.70 ng/u1

response 8566

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.30	36.17
69.00	20.30	27.47#
0.00	0.00	0.00

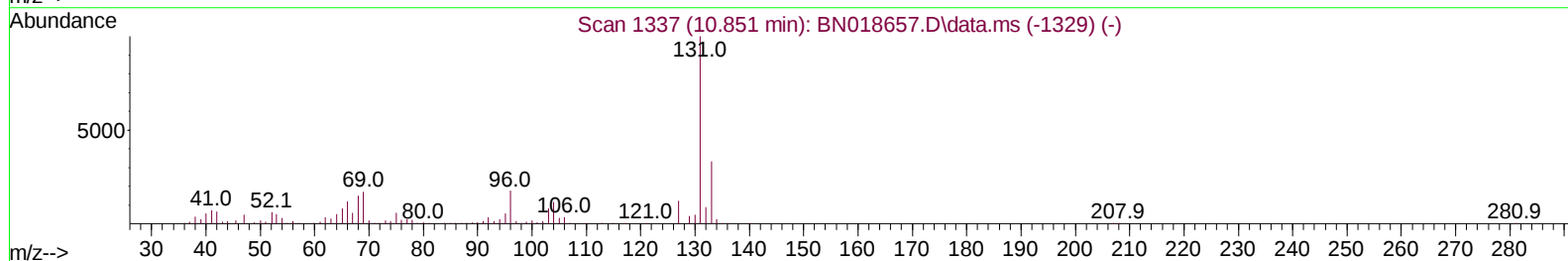
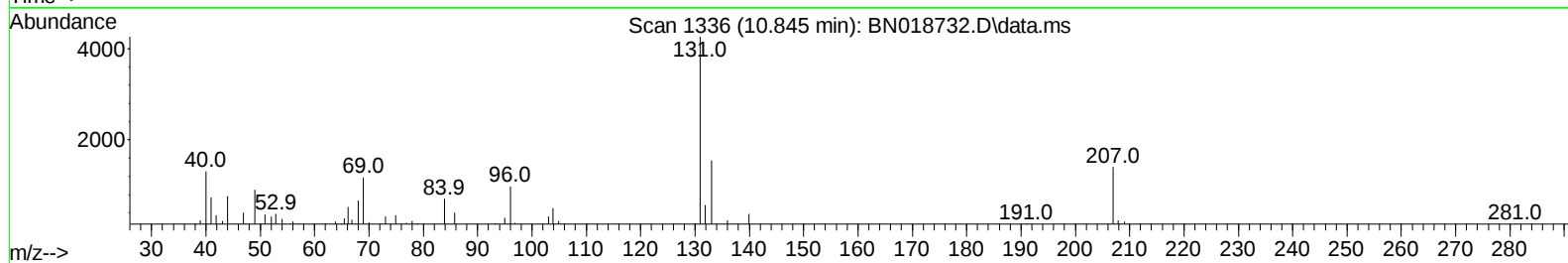
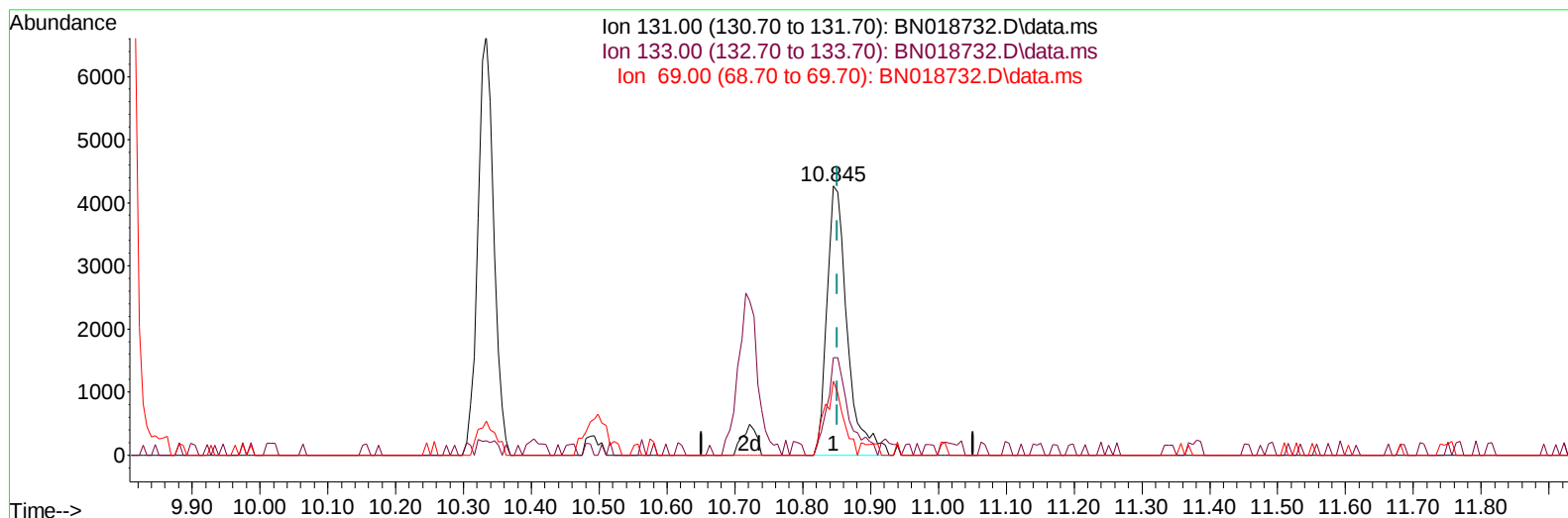
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TIC: BN018732.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.845min (-0.006) 0.73 ng/ul m

response 8877

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	33.30	36.17
69.00	20.30	27.47#
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	7.910	152	149402	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	582349	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.545	164	335623	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	643887	20.000	ng/ul	0.00
79) Chrysene-d12	21.462	240	625151	20.000	ng/ul	0.00
88) Perylene-d12	23.868	264	689718	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	23336	5.887	ng/uL	0.00
4) Pyridine-d5	3.710	84	43784	4.086	ng/ul	0.00
7) Phenol-d5	7.063	99	96611	7.841	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.228	67	298643	39.051	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	296536	30.556	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	173970	18.488	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	168988	37.764	ng/ul	0.00
24) 2-Nitrophenol-d4	9.792	143	176053	36.735	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.334	165	298705	32.655	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	8877m	0.726	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	962584	39.619	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	1101779	35.002	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	26632	6.308	ng/ul	0.00
60) Fluorene-d10	15.533	176	869612	41.457	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	121679	32.029	ng/ul	0.00
73) Anthracene-d10	17.380	188	1382464	44.871	ng/ul	0.00
81) Pyrene-d10	19.674	212	1592235	43.514	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.709	264	1591405	45.243	ng/ul	0.00
Target Compounds						
2) 1,4-Dioxane	3.322	88	13731	3.278	ng/uL	Qvalue 98

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

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Manual IntegrationsAPPROVED

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