

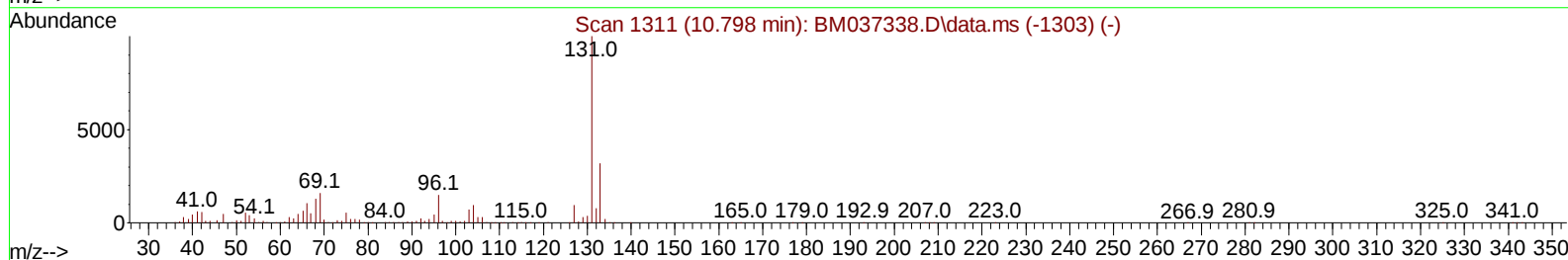
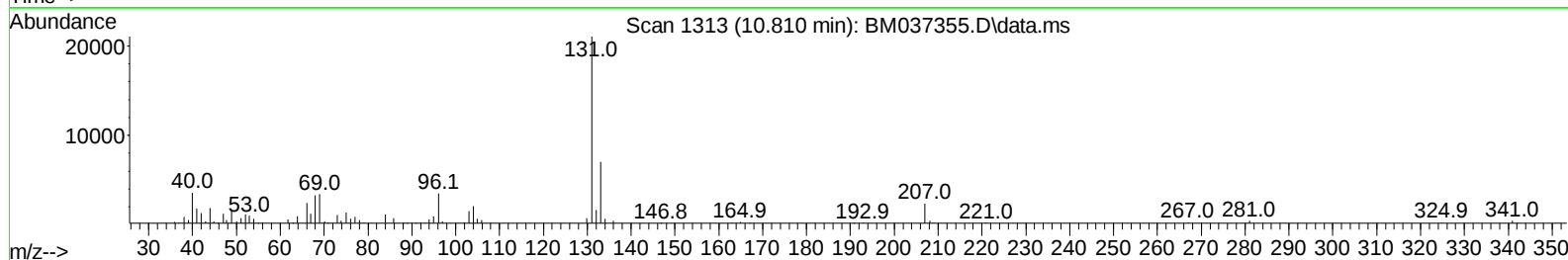
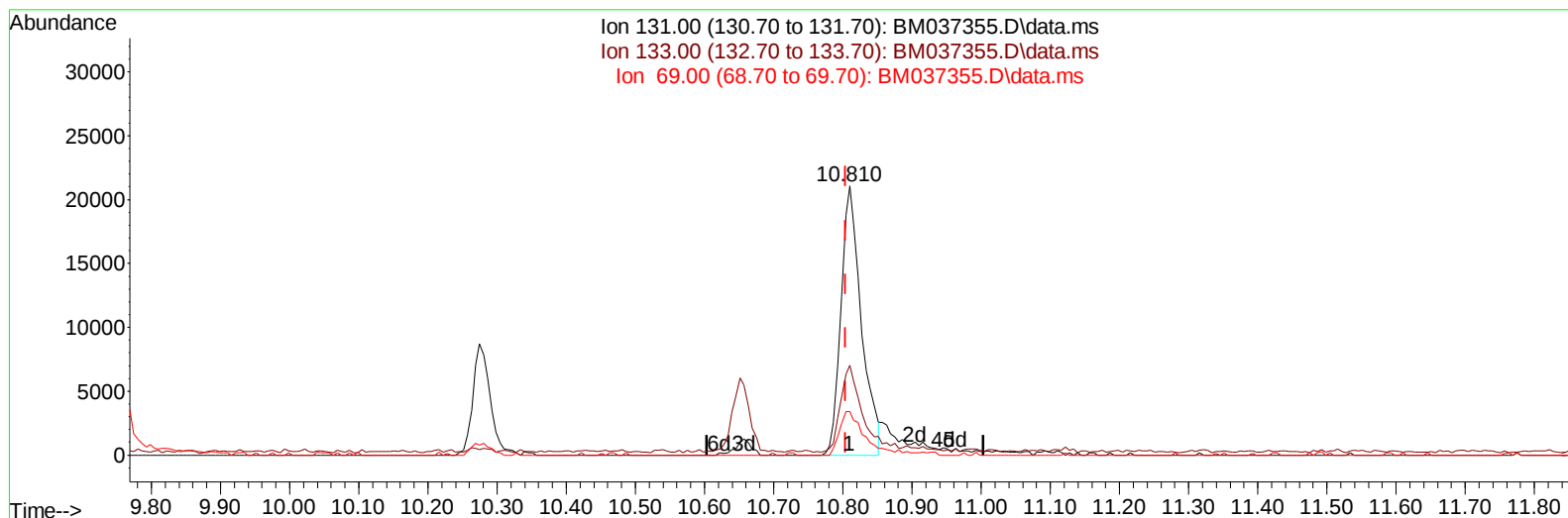
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
Data File : BM037355.D
Acq On : 03 Nov 2022 20:16
Operator : CG/JU
Sample : N5276-01
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4W2

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 11/04/2022
Supervised By :Jagrut Upadhyay 11/07/2022

Quant Time: Nov 04 01:06:09 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 02 03:33:37 2022
Response via : Initial Calibration



TIC: BM037355.D\data.ms

(31) 4-Chloroaniline-d4 (S)**10.810min (+ 0.006) 1.40 ng/u1****response 43163**

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.44
69.00	16.10	16.11
0.00	0.00	0.00

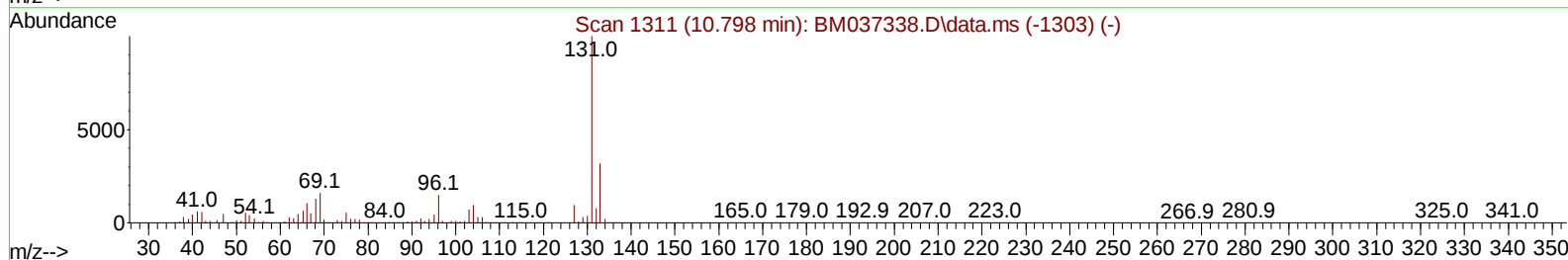
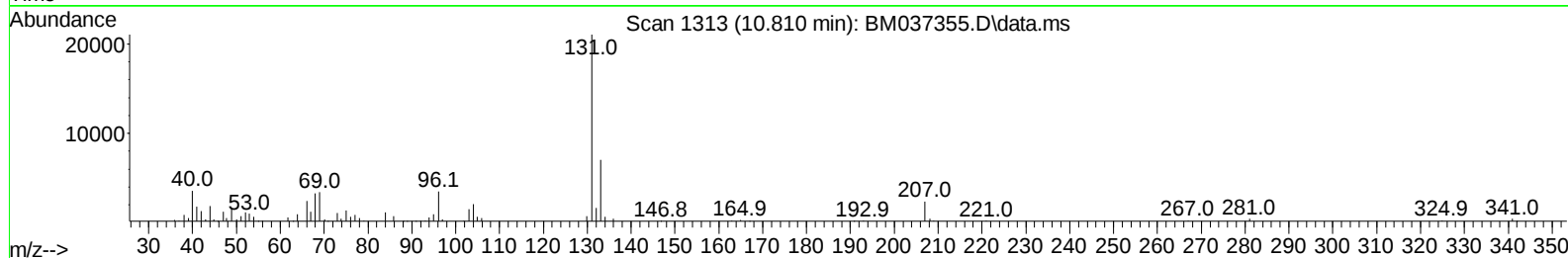
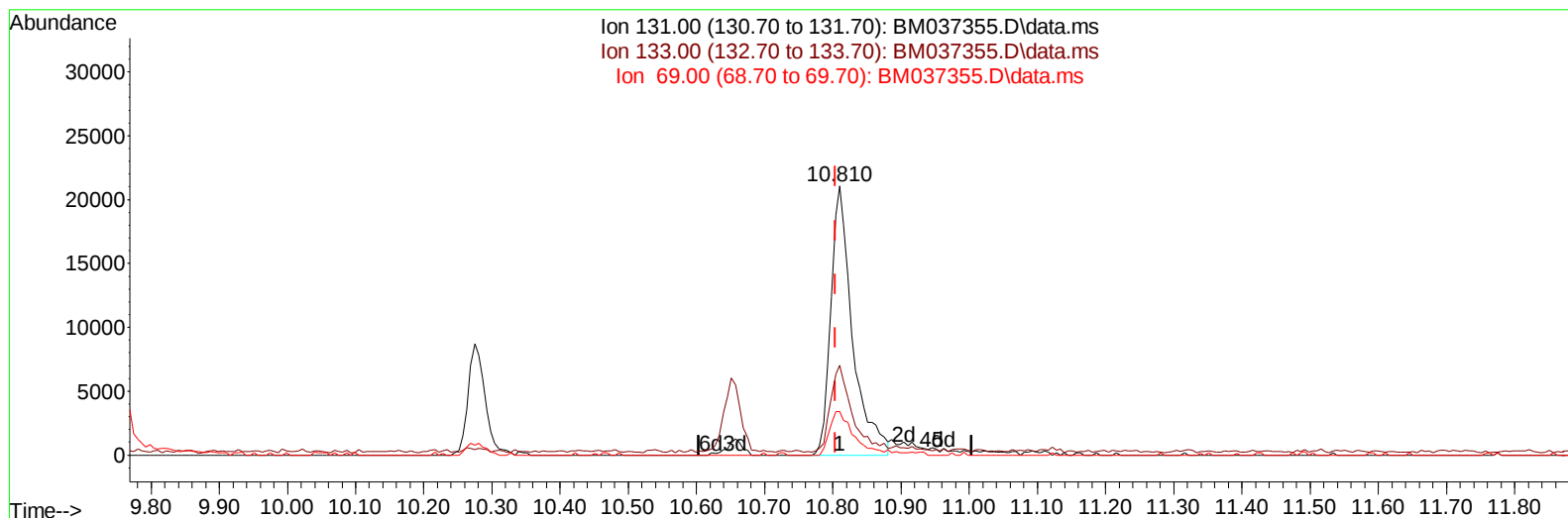
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TIC: BM037355.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.810min (+ 0.006) 1.51 ng/ul m

response 46345

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	33.44
69.00	16.10	16.11
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	7.851	152	295691	20.000	ng/ul	0.00
20) Naphthalene-d8	10.651	136	1342042	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.486	164	876935	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.227	188	1824322	20.000	ng/ul	0.00
79) Chrysene-d12	21.398	240	1700041	20.000	ng/ul	0.00
88) Perylene-d12	23.744	264	1485764	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	31664	4.720	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.022	99	184728	7.179	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	497828	31.526	ng/ul	0.00
11) 2-Chlorophenol-d4	7.381	132	492239	25.036	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	330127	15.710	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	319178	31.839	ng/ul	0.00
24) 2-Nitrophenol-d4	9.739	143	319192	29.230	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	540889	26.866	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	46345m	1.506	ng/ul	0.00
46) Dimethylphthalate-d6	13.904	166	2136339	35.234	ng/ul	0.00
49) Acenaphthylene-d8	14.180	160	2396553	32.877	ng/ul	0.00
54) 4-Nitrophenol-d4	14.710	143	66590	5.494	ng/ul	0.00
60) Fluorene-d10	15.474	176	2016267	37.290	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	311825	26.986	ng/ul	0.00
73) Anthracene-d10	17.327	188	3378939	40.217	ng/ul	0.00
81) Pyrene-d10	19.615	212	3775928	42.504	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.592	264	3092987	40.442	ng/ul	-0.01

Target Compounds Qvalue

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

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