

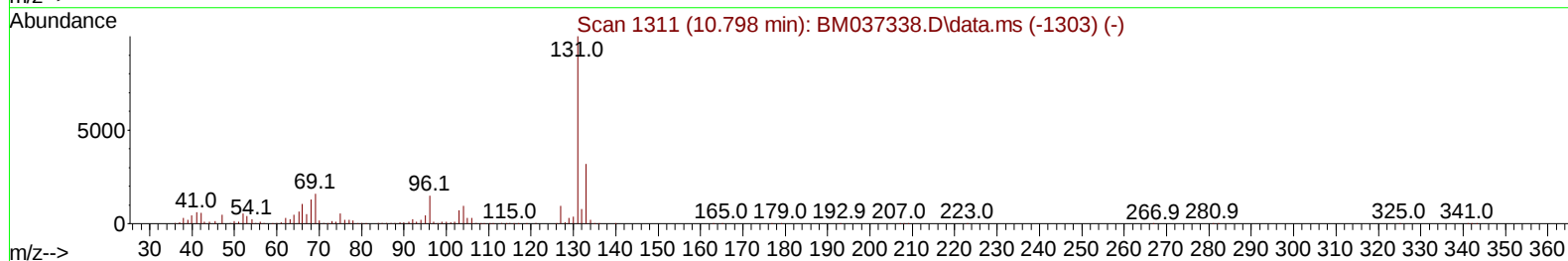
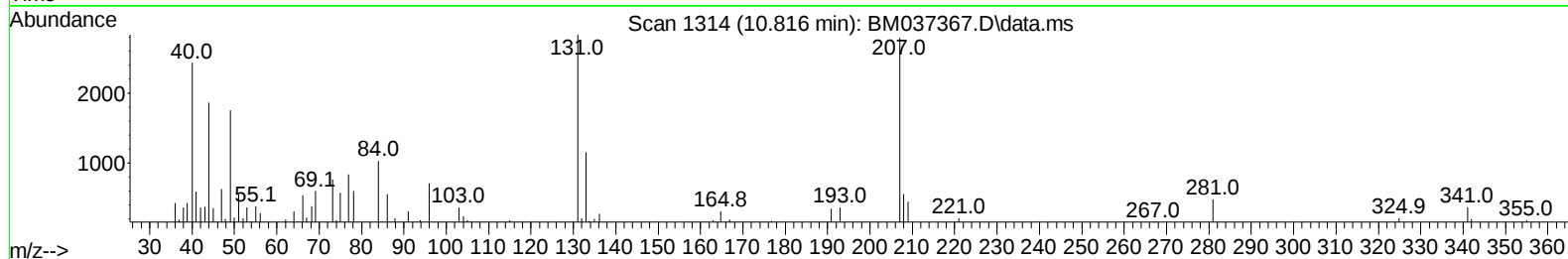
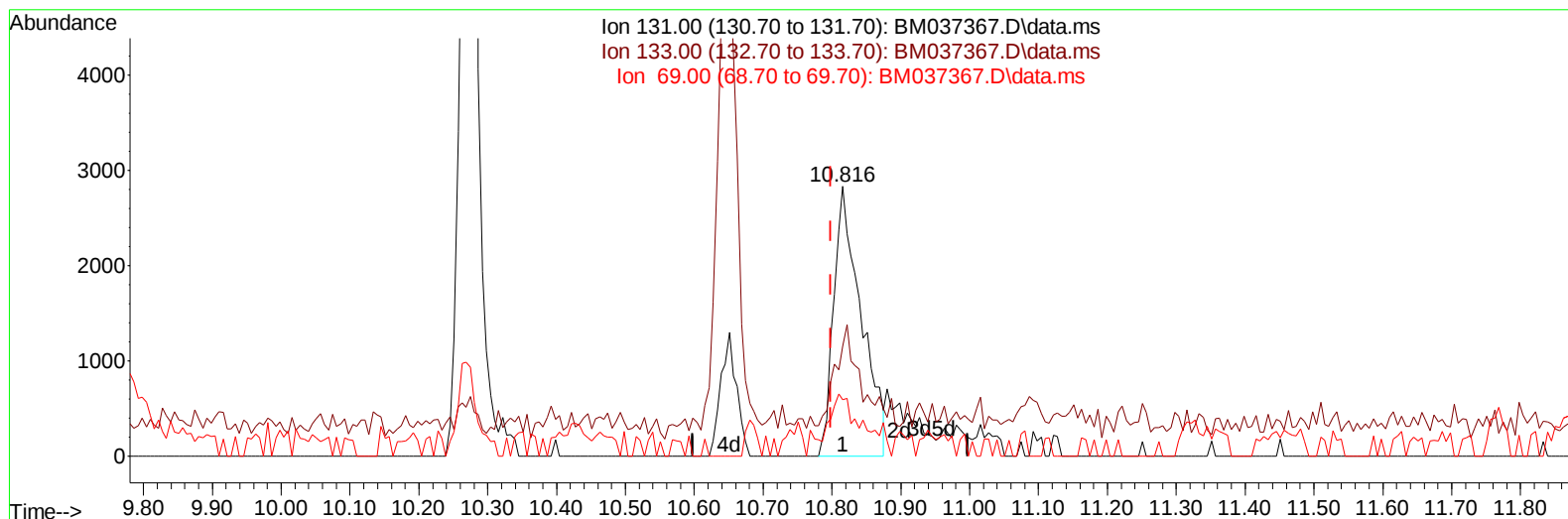
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM110322\
 Data File : BM037367.D
 Acq On : 04 Nov 2022 03:55
 Operator : CG/JU
 Sample : N5276-05
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4W5

Manual IntegrationsAPPROVED

Quant Time: Nov 04 04:38:19 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM110122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 04 02:39:09 2022
 Response via : Initial Calibration

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



TIC: BM037367.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.816min (+ 0.018) 0.24 ng/ul

response 7803

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	40.68#
69.00	16.10	21.01#
0.00	0.00	0.00

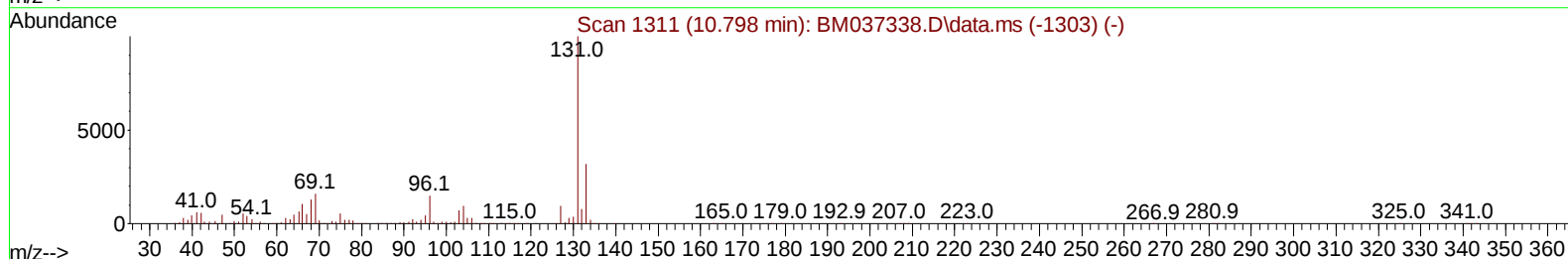
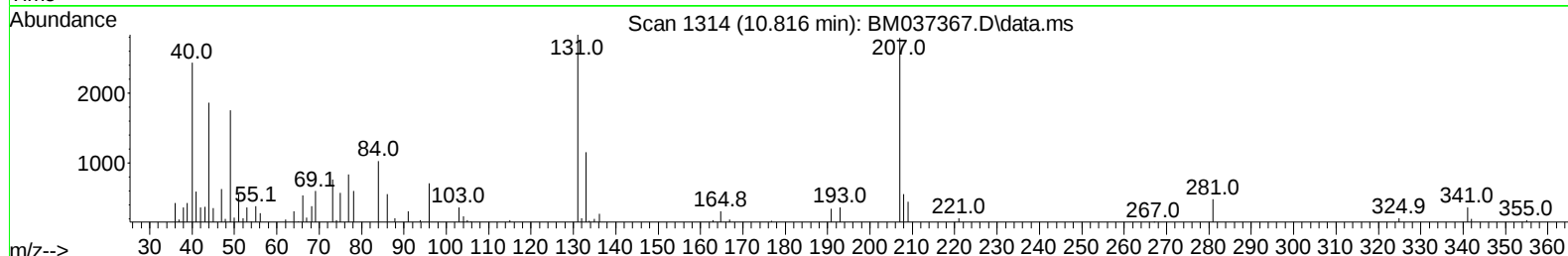
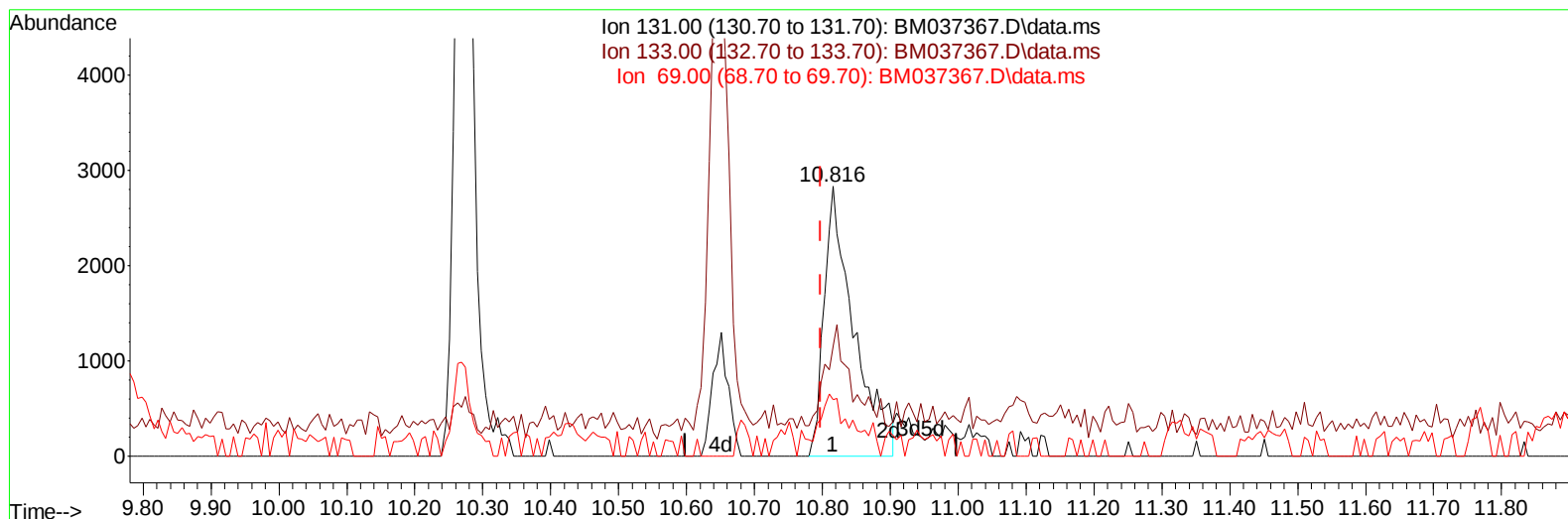
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TIC: BM037367.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.816min (+ 0.018) 0.27 ng/ul m

response 8725

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	40.68#
69.00	16.10	21.01#
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	7.845	152	318944	20.000	ng/ul	0.00
20) Naphthalene-d8	10.645	136	1430909	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.480	164	925382	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.221	188	1909467	20.000	ng/ul	0.00
79) Chrysene-d12	21.391	240	1732858	20.000	ng/ul	0.00
88) Perylene-d12	23.738	264	1471250	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	27011	3.733	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.022	99	179368	6.462	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	463094	27.188	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	474262	22.363	ng/ul	0.00
15) 4-Methylphenol-d8	8.563	113	333565	14.717	ng/ul	0.00
21) Nitrobenzene-d5	9.016	128	295641	27.659	ng/ul	0.00
24) 2-Nitrophenol-d4	9.733	143	302282	25.962	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.275	165	548517	25.553	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	8725m	0.266	ng/ul	0.02
46) Dimethylphthalate-d6	13.898	166	1909261	29.840	ng/ul	0.00
49) Acenaphthylene-d8	14.174	160	2282943	29.679	ng/ul	0.00
54) 4-Nitrophenol-d4	14.704	143	48142	3.764	ng/ul	0.00
60) Fluorene-d10	15.474	176	1831511	32.099	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.598	200	265243	21.931	ng/ul	0.00
73) Anthracene-d10	17.321	188	2912331	33.117	ng/ul	0.00
81) Pyrene-d10	19.609	212	3215838	35.514	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.585	264	2522736	33.311	ng/ul	0.00

Target Compounds Qvalue

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

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