

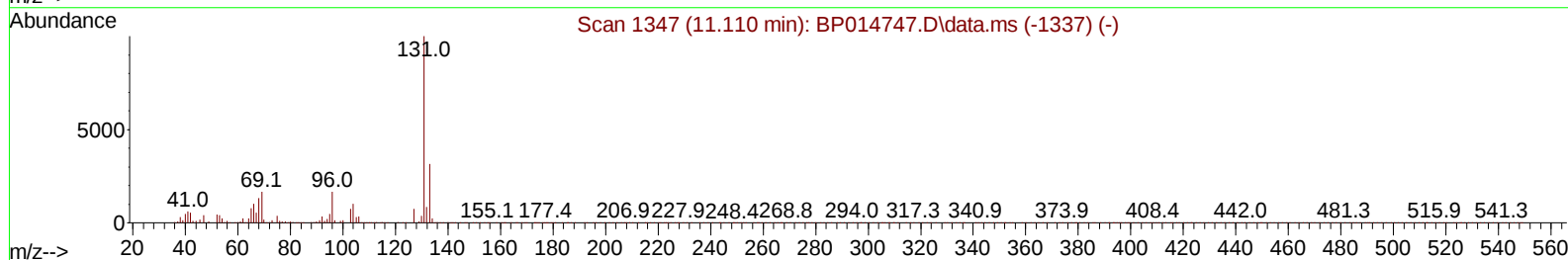
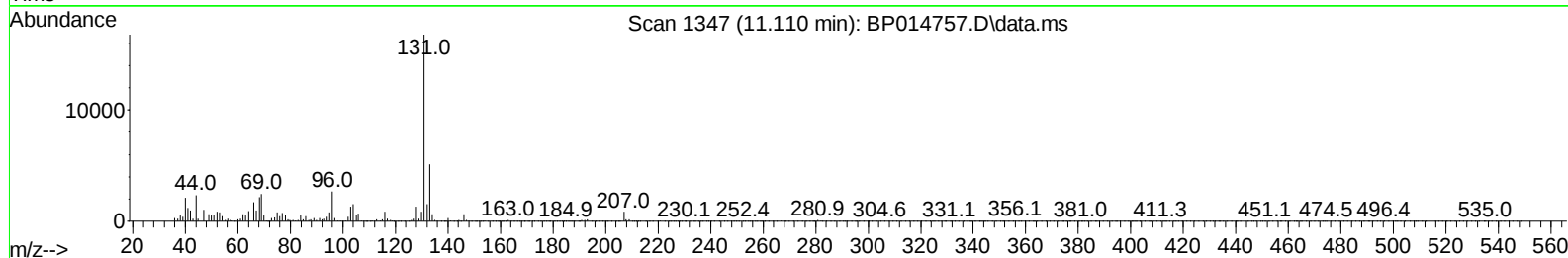
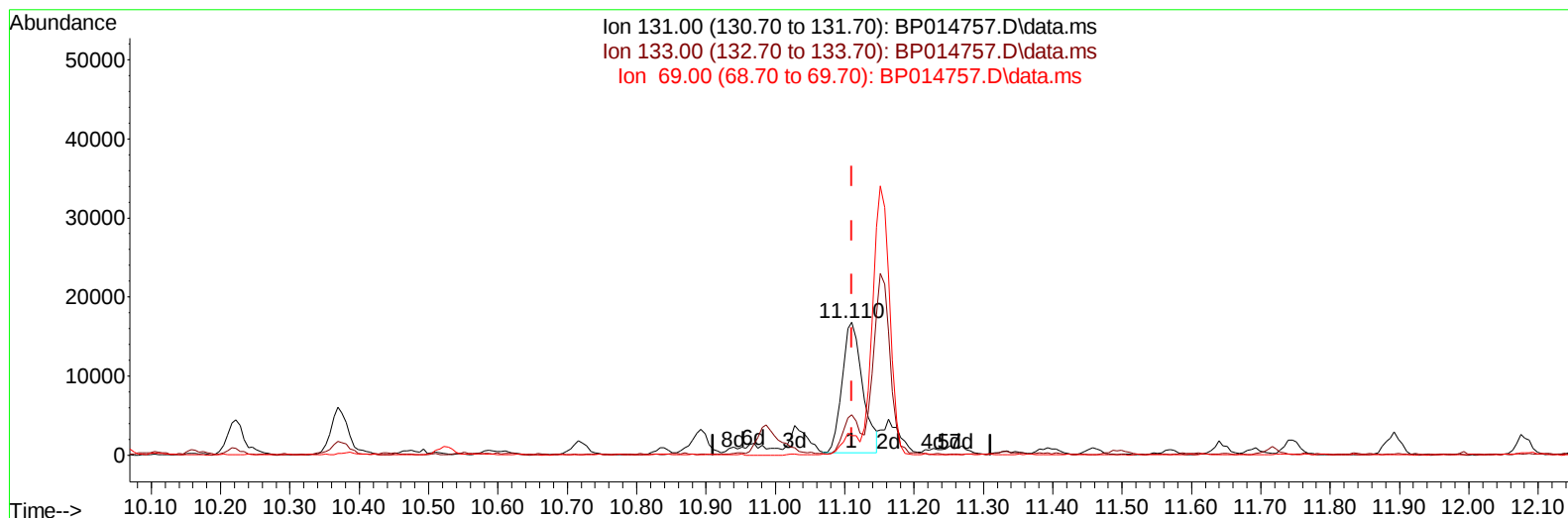
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042023\
 Data File : BP014757.D
 Acq On : 20 Apr 2023 15:43
 Operator : CG/JU
 Sample : 02296-08DL 10X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCBL0DL

Manual IntegrationsAPPROVED

Quant Time: Apr 21 03:55:22 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 20 02:32:50 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/21/2023
 Supervised By :Jagrut Upadhyay 04/21/2023



TIC: BP014757.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.110min (-0.000) 1.61 ng/u1

response 33637

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.90	30.51
69.00	17.90	14.71
0.00	0.00	0.00

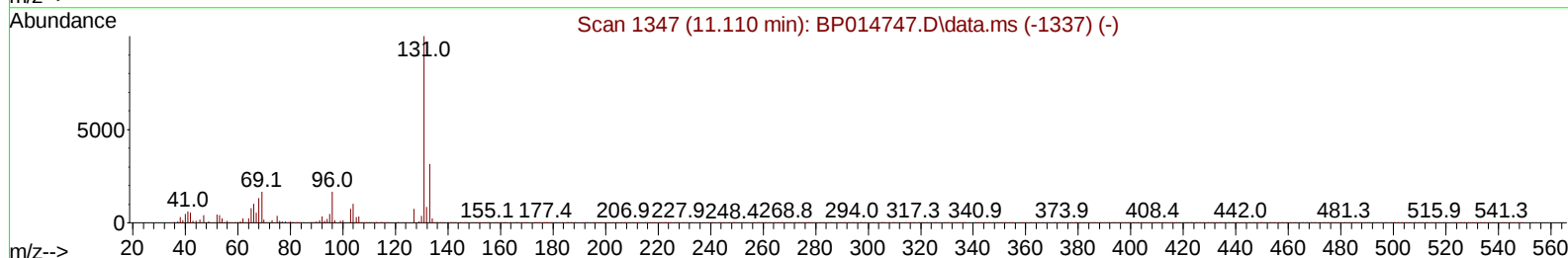
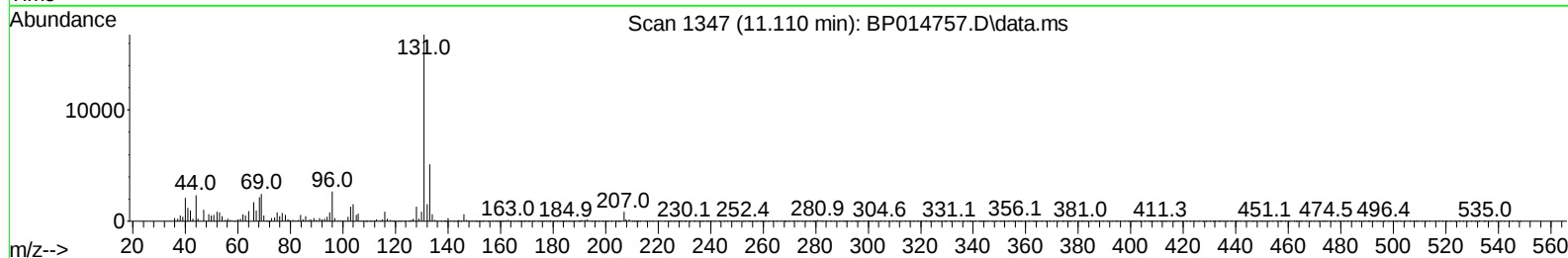
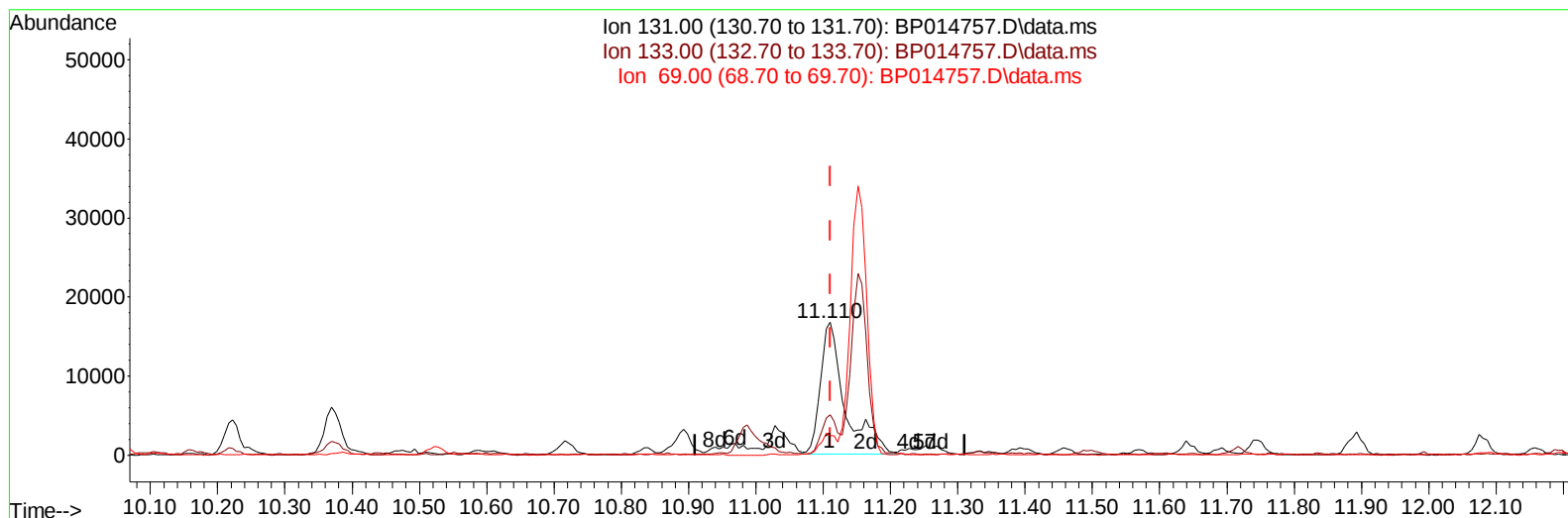
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Manual IntegrationsAPPROVED

Quant Time: Apr 20 22:40:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP041823.M
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TIC: BP014757.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.110min (-0.000) 2.03 ng/ul m

response 42547

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.90	30.51
69.00	17.90	14.71
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	8.152	152	221595	20.000	ng/ul	0.00
20) Naphthalene-d8	10.987	136	893176	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.781	164	528839	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.533	188	1145826	20.000	ng/ul	0.00
79) Chrysene-d12	21.604	240	1072360	20.000	ng/ul	0.00
88) Perylene-d12	24.180	264	1215322	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.487	96	2473	0.444	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.287	99	8694	0.454	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.463	67	33563	2.840	ng/ul	0.00
11) 2-Chlorophenol-d4	7.675	132	25740	1.809	ng/ul	0.00
15) 4-Methylphenol-d8	8.851	113	16334	1.094	ng/ul	0.00
21) Nitrobenzene-d5	9.322	128	15181	2.712	ng/ul	0.00
24) 2-Nitrophenol-d4	10.051	143	12201	2.414	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.587	165	28454	2.073	ng/ul	0.00
31) 4-Chloroaniline-d4	11.110	131	42547m	2.031	ng/ul	0.00
46) Dimethylphthalate-d6	14.181	166	125874	3.006	ng/ul	0.00
49) Acenaphthylene-d8	14.475	160	125318	2.582	ng/ul	0.00
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.769	176	106361	2.960	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	0.000	200	0d	0.000	ng/ul	
73) Anthracene-d10	17.628	188	172969	3.125	ng/ul	0.00
81) Pyrene-d10	19.845	212	191192	3.112	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.004	264	179167	2.844	ng/ul	-0.01
Target Compounds						
					Qvalue	
30) Naphthalene	11.046	128	17802536	355.643	ng/ul	98
36) 2-Methylnaphthalene	12.622	142	891161	26.667	ng/ul	100
37) 1-Methylnaphthalene	12.840	142	1704872	51.461	ng/ul	99
43) 1,1'-Biphenyl	13.622	154	640301	14.932	ng/ul	98
52) Acenaphthene	14.845	153	1690090	47.130	ng/ul	99
56) Dibenzofuran	15.175	168	2029889	40.219	ng/ul	99
61) Fluorene	15.828	166	799679	19.411	ng/ul	99
72) Phenanthrene	17.575	178	1256593	19.459	ng/ul	99
77) Carbazole	17.922	167	1429519	24.873	ng/ul	99

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

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