

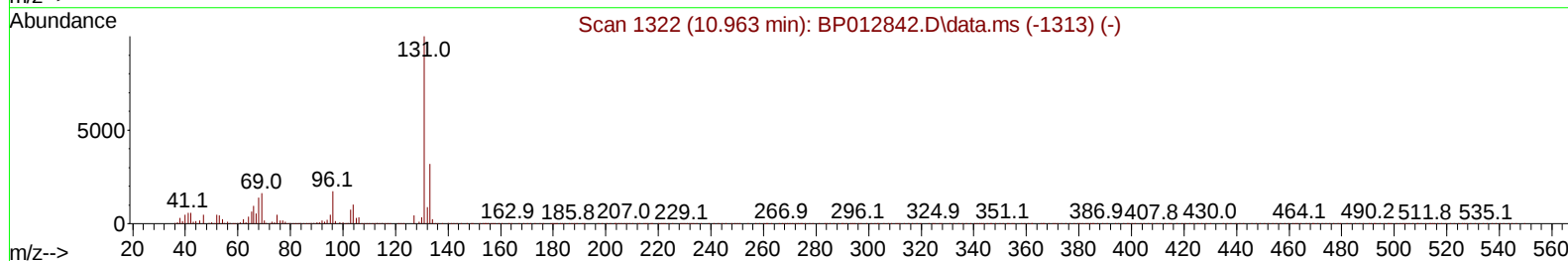
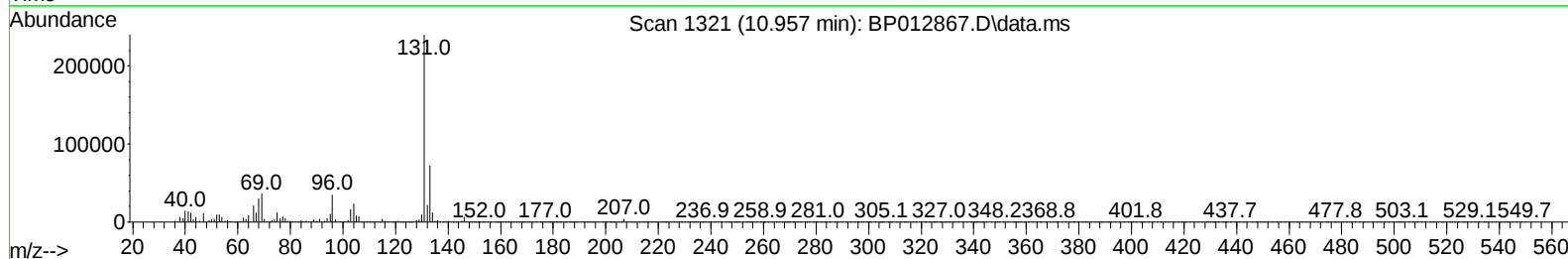
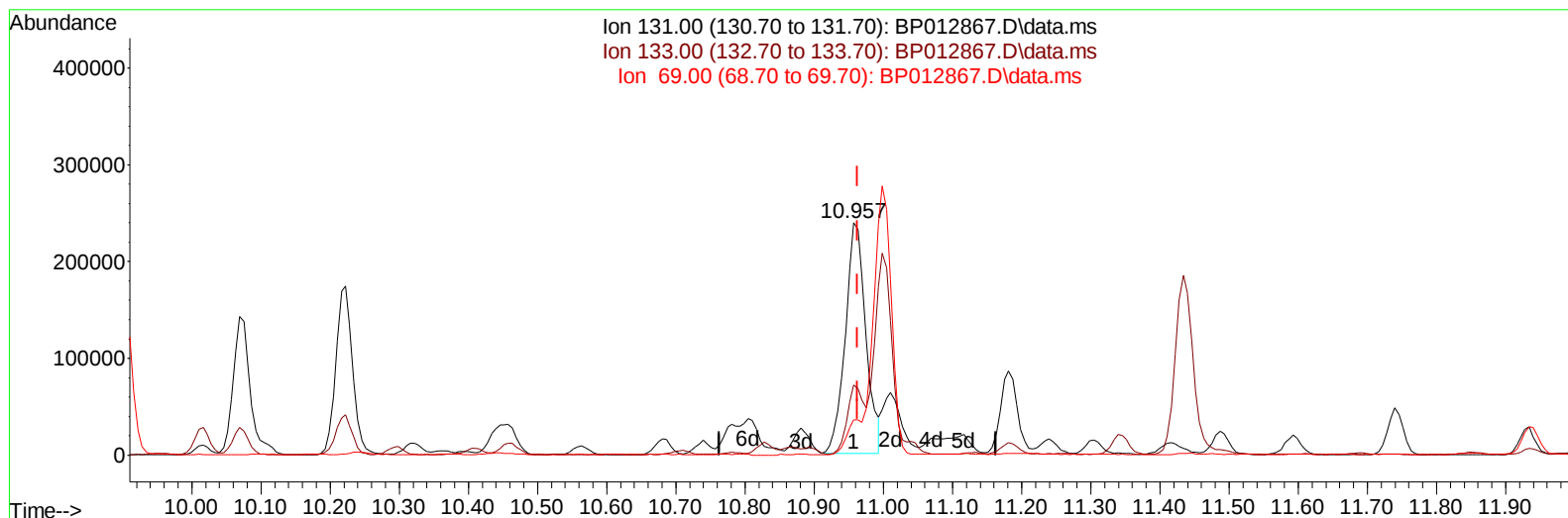
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112922\
 Data File : BP012867.D
 Acq On : 29 Nov 2022 23:44
 Operator : CG/JU
 Sample : N5725-07
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BHB63

Manual IntegrationsAPPROVED

Quant Time: Nov 30 00:50:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 29 22:53:34 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/30/2022
 Supervised By :mohammad ahmed 11/30/2022



TIC: BP012867.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.957min (-0.006) 8.67 ng/u1

response 489983

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.40	30.24
69.00	15.70	15.15
0.00	0.00	0.00

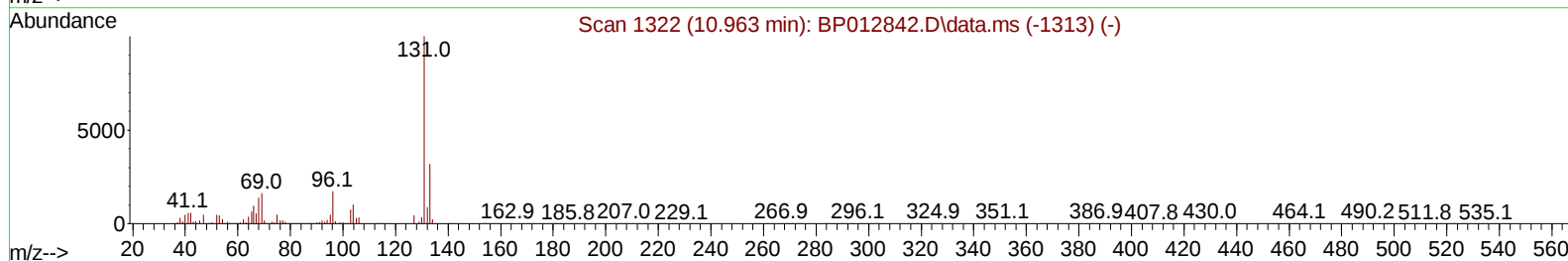
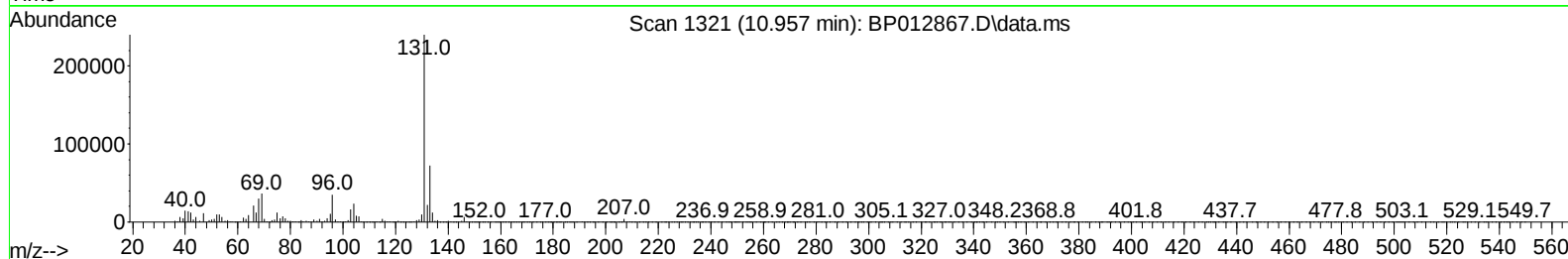
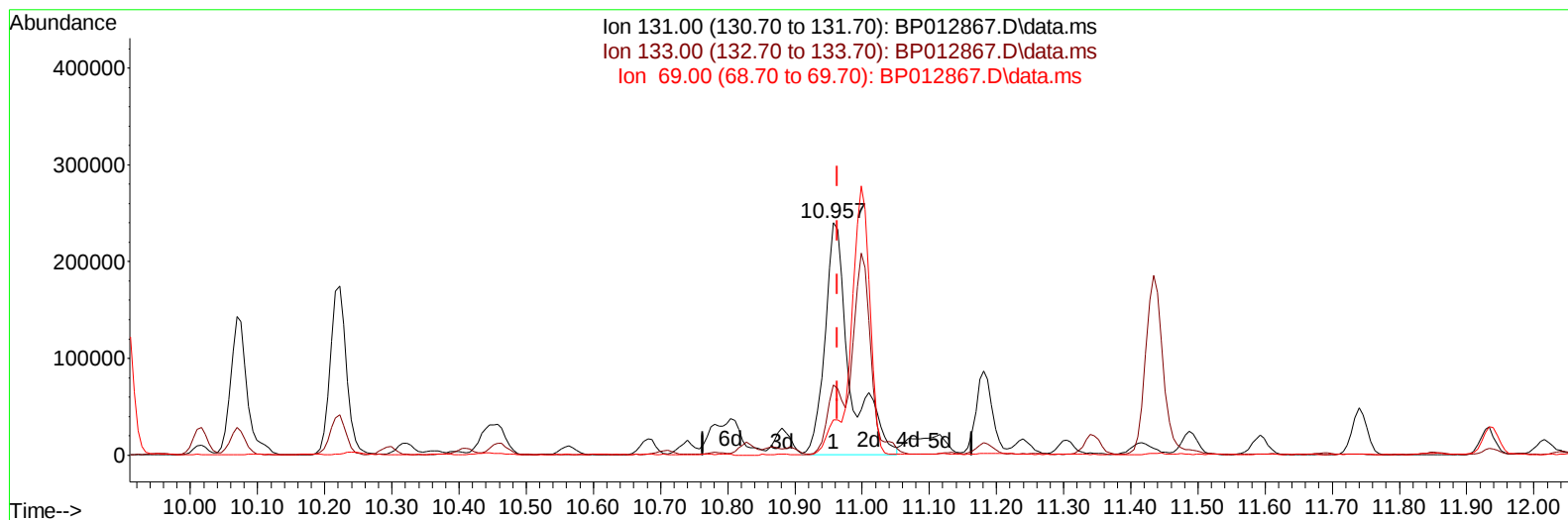
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TIC: BP012867.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.957min (-0.006) 10.99 ng/ul m

response 621278

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.40	30.24
69.00	15.70	15.15
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	8.010	152	619068	20.000	ng/ul	0.00
20) Naphthalene-d8	10.828	136	2723091	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.645	164	1692235	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.398	188	2898987	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	1929035	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264	2160926	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.399	96	67728	4.688	ng/uL	0.00
4) Pyridine-d5	3.822	84	357973	8.280	ng/ul	0.00
7) Phenol-d5	7.163	99	395366	7.772	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.334	67	962398	30.702	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	994895	25.666	ng/ul	0.00
15) 4-Methylphenol-d8	8.710	113	699333	17.062	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	623807	37.428	ng/ul	0.00
24) 2-Nitrophenol-d4	9.904	143	607120	34.306	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.440	165	1211112	29.330	ng/ul	0.00
31) 4-Chloroaniline-d4	10.957	131	621278m	10.987	ng/ul	0.00
46) Dimethylphthalate-d6	14.051	166	3818956	32.786	ng/ul	0.00
49) Acenaphthylene-d8	14.340	160	4405736	32.417	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	84451	4.067	ng/ul	0.00
60) Fluorene-d10	15.639	176	3327665	32.621	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.745	200	415697	30.002	ng/ul	0.00
73) Anthracene-d10	17.498	188	4500992	35.271	ng/ul	0.00
81) Pyrene-d10	19.727	212	4048723	36.121	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.827	264	3717311	35.022	ng/ul	0.00
Target Compounds						Qvalue
30) Naphthalene	10.881	128	2491967	17.476	ng/ul	99
37) 1-Methylnaphthalene	12.698	142	2067240	21.007	ng/ul	99
43) 1,1'-Biphenyl	13.481	154	2544104	20.638	ng/ul	100
52) Acenaphthene	14.710	153	5719417	54.666	ng/ul	99
56) Dibenzofuran	15.045	168	5394191	37.405	ng/ul	99
61) Fluorene	15.692	166	161697	1.398	ng/ul	96
72) Phenanthrene	17.445	178	581540	3.764	ng/ul	100

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

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