

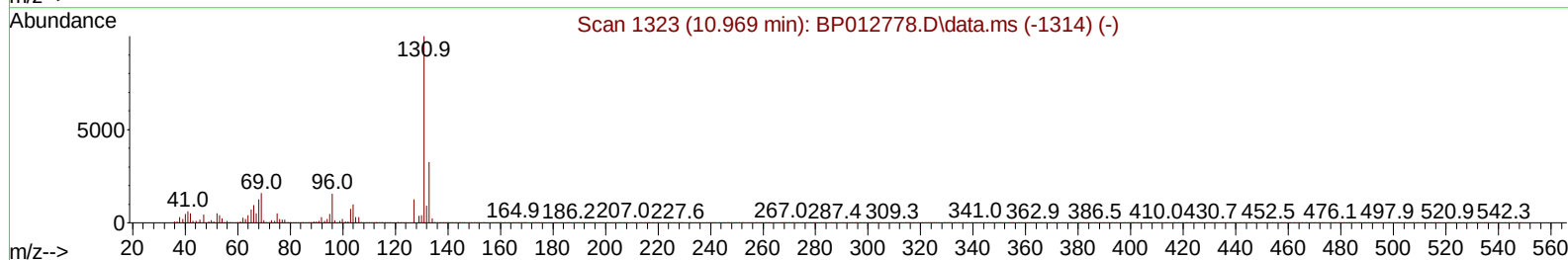
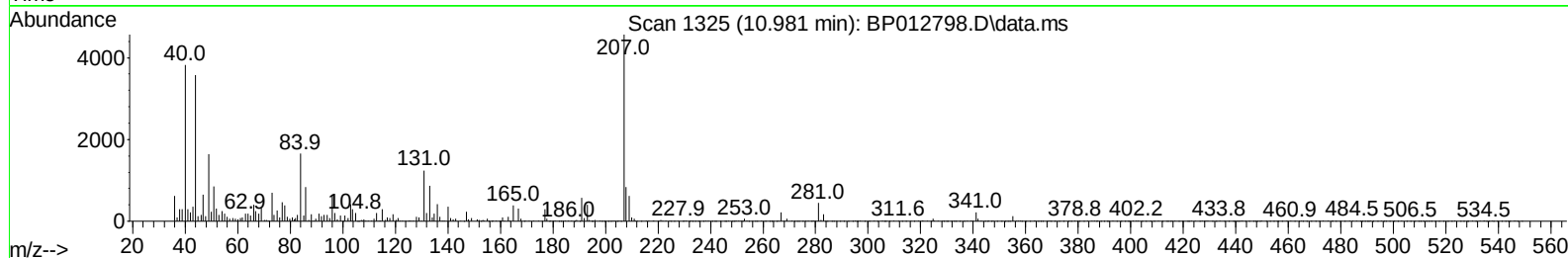
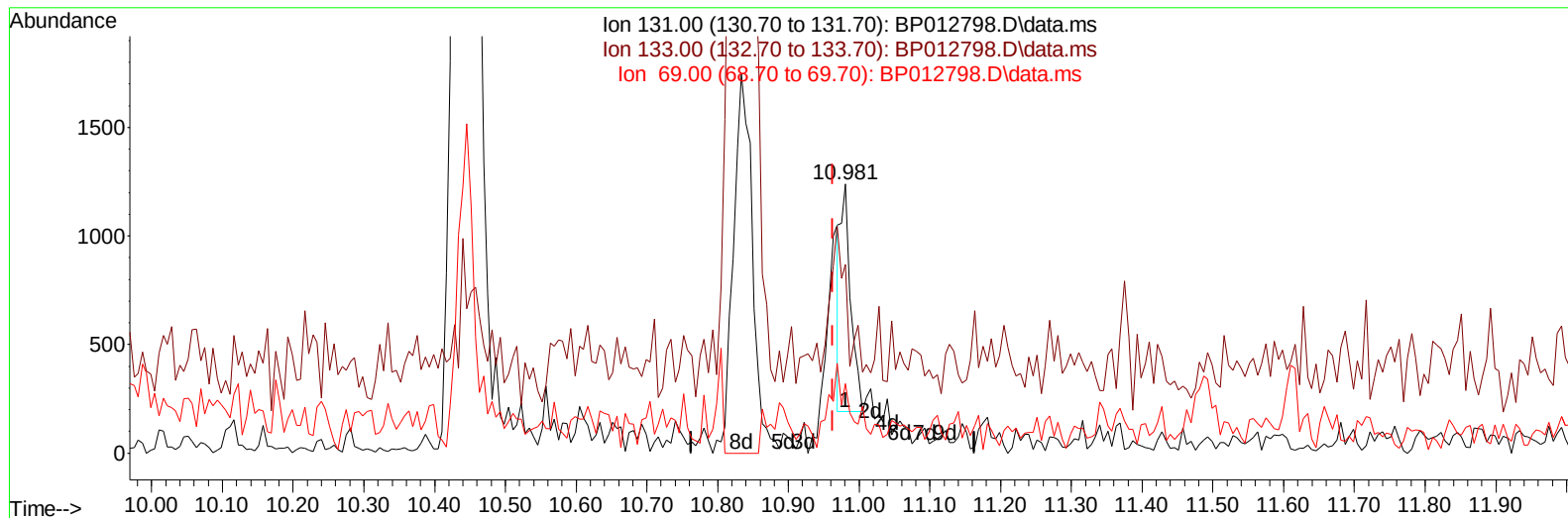
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP112522\
 Data File : BP012798.D
 Acq On : 27 Nov 2022 09:29
 Operator : CG/JU
 Sample : PB149206BL
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK206

Manual IntegrationsAPPROVED

Quant Time: Dec 03 04:39:42 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 11/28/2022
 Supervised By :mohammad ahmed 11/30/2022



TIC: BP012798.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.981min (+ 0.018) 0.02 ng/u1

response 1039

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.40	70.14#
69.00	15.70	25.83#
0.00	0.00	0.00

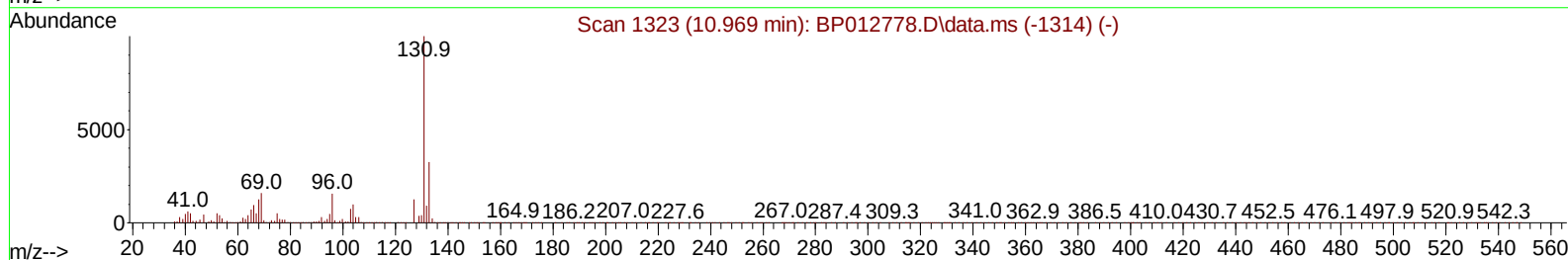
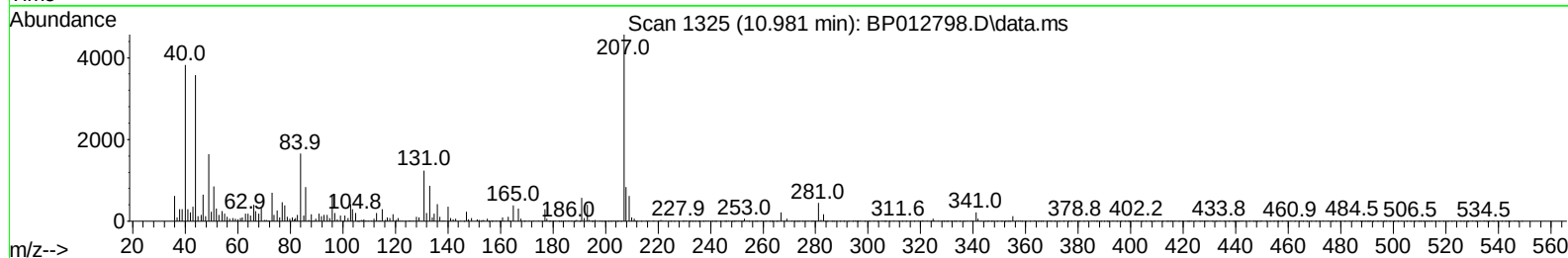
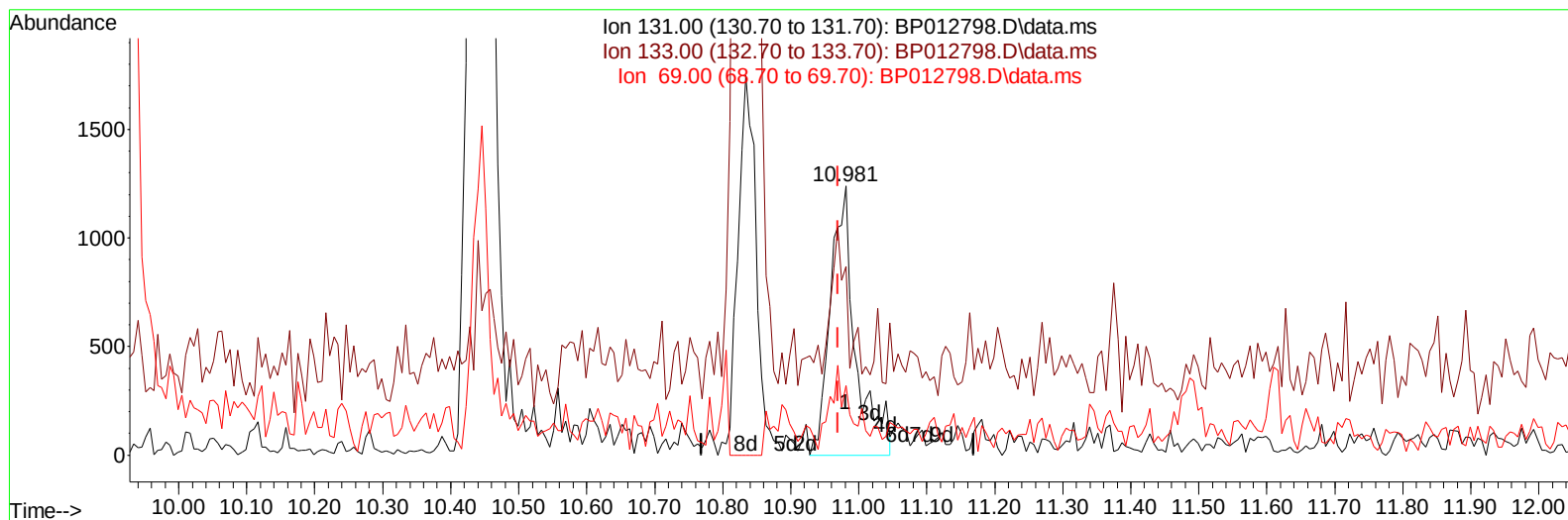
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 Acq On : 27 Nov 2022 09:29
 Operator : CG/JU
 Sample : PB149206BL
 Misc :
 ALS Vial : 73 Sample Multiplier: 1

Instrument :
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 ClientSampleId :
 SBLK206

Manual IntegrationsAPPROVED

Quant Time: Nov 28 00:02:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP112522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Nov 27 22:55:08 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/28/2022
 Supervised By :mohammad ahmed 11/30/2022



TIC: BP012798.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.981min (+ 0.012) 0.07 ng/ul m

response 3185

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.40	70.14#
69.00	15.70	25.83#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.016	152	500721	20.000	ng/ul	0.00
20) Naphthalene-d8	10.834	136	2270895	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1573404	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.410	188	3508153	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	3439566	20.000	ng/ul	0.00
88) Perylene-d12	24.010	264	3405327	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	61477	5.261	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.164	99	1092950	26.563	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.340	67	717623	28.304	ng/ul	0.00
11) 2-Chlorophenol-d4	7.540	132	884419	28.209	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113	900427	27.161	ng/ul	0.00
21) Nitrobenzene-d5	9.181	128	430724	30.989	ng/ul	0.00
24) 2-Nitrophenol-d4	9.905	143	479793	32.510	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.446	165	920628	26.735	ng/ul	0.00
31) 4-Chloroaniline-d4	10.981	131	3185m	0.068	ng/ul	0.01
46) Dimethylphthalate-d6	14.057	166	3355883	30.987	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	3575009	28.291	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143	489538	25.354	ng/ul	0.00
60) Fluorene-d10	15.645	176	2875670	30.319	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	386188	23.032	ng/ul	0.00
73) Anthracene-d10	17.510	188	4682274	30.321	ng/ul	0.00
81) Pyrene-d10	19.733	212	5468313	27.361	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.845	264	5048286	30.181	ng/ul	0.00

Target Compounds Qvalue

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

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