

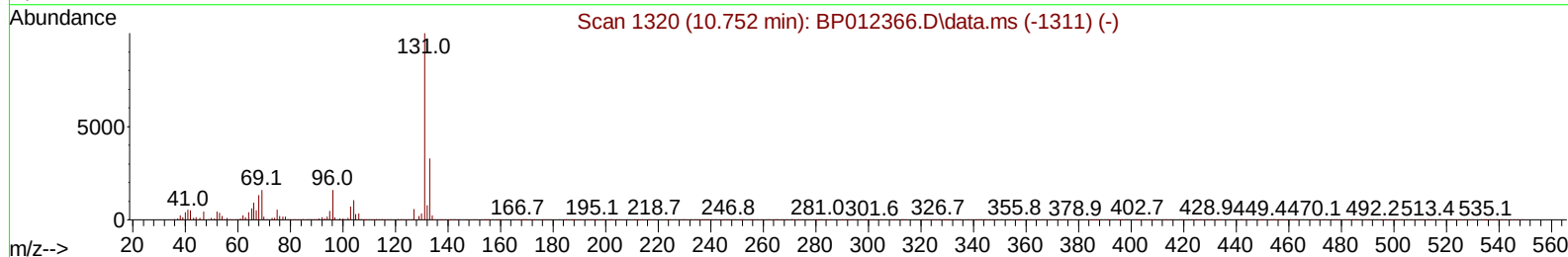
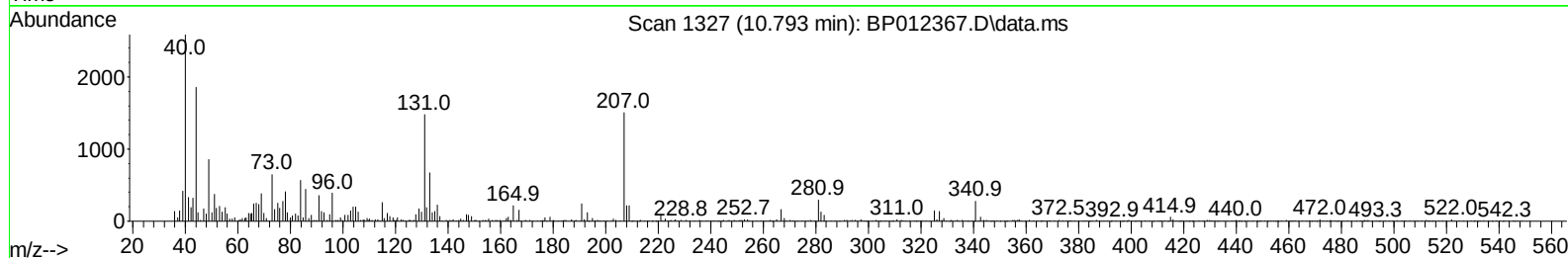
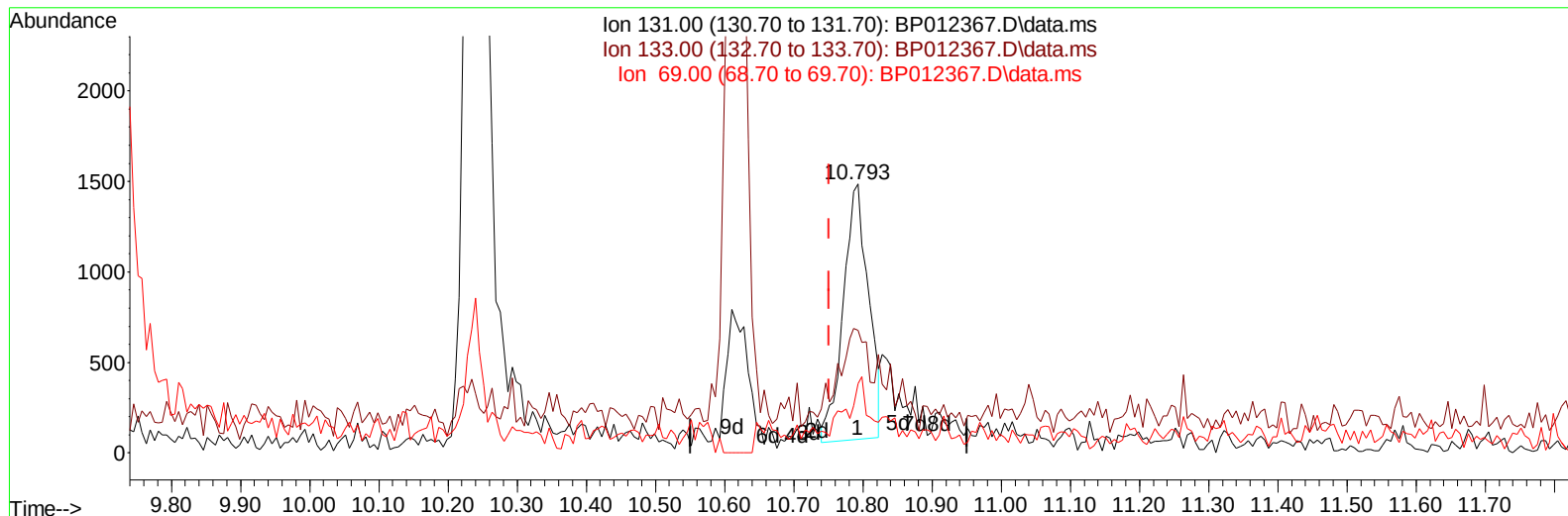
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102722\
Data File : BP012367.D
Acq On : 28 Oct 2022 08:58
Operator : CG/JU
Sample : PB148502BL
Misc :
ALS Vial : 30 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK502

Manual IntegrationsAPPROVED

Quant Time: Oct 28 23:18:53 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 28 06:05:59 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/31/2022
Supervised By :mohammad ahmed 10/31/2022



TIC: BP012367.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.793min (+ 0.041) 0.15 ng/u1

response 3559

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	45.72#
69.00	16.50	25.79#
0.00	0.00	0.00

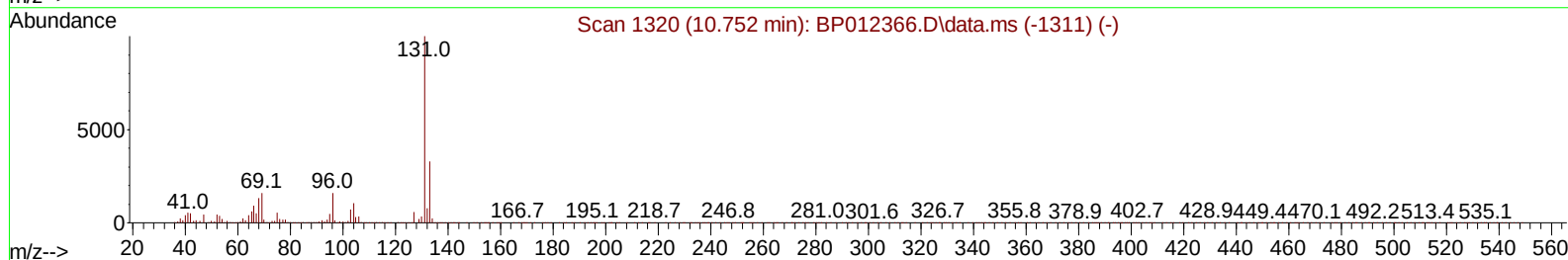
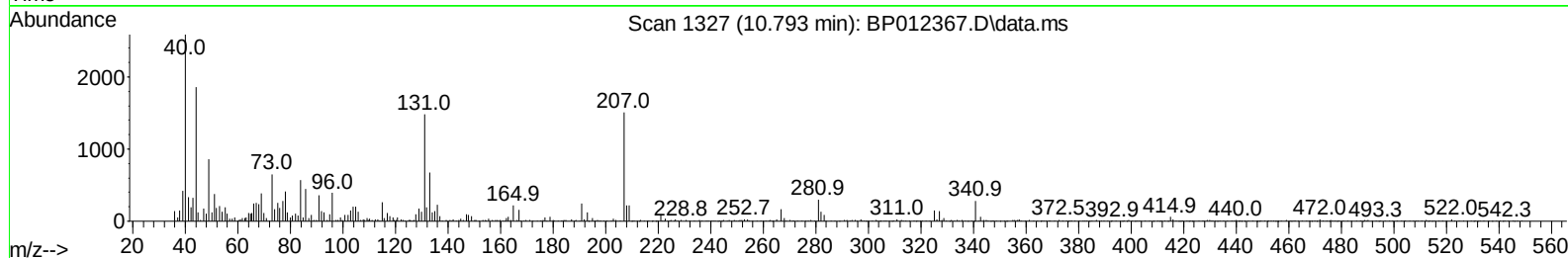
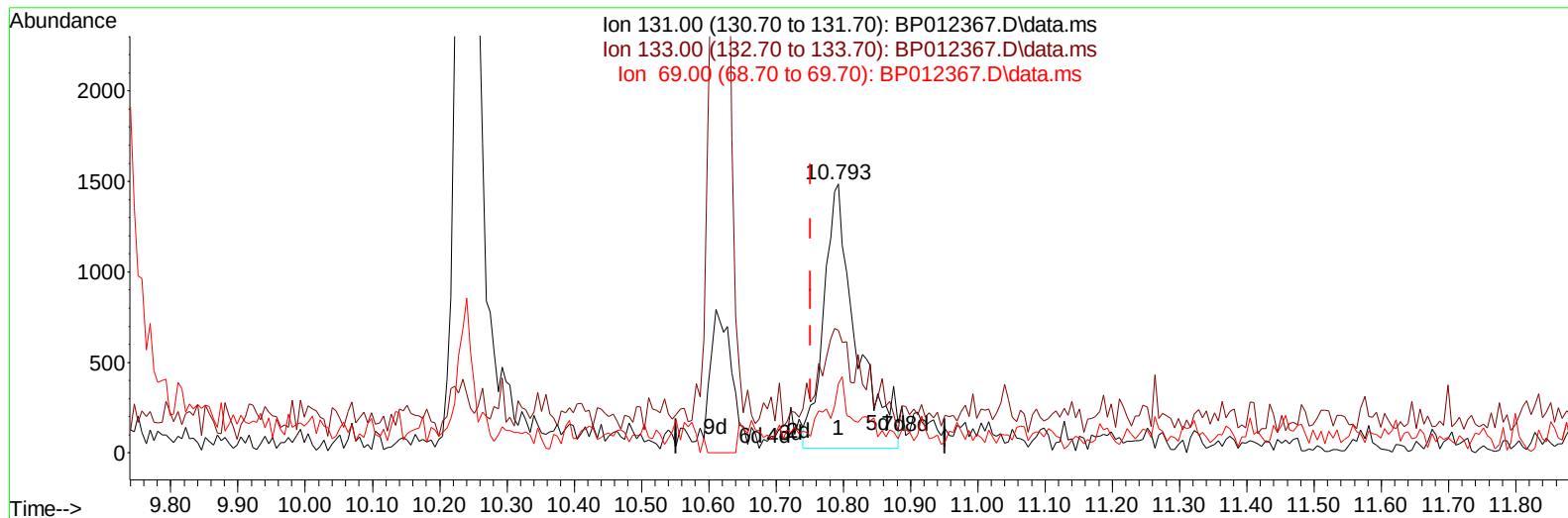
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TIC: BP012367.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.793min (+ 0.041) 0.21 ng/ul m

response 4892

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	45.72#
69.00	16.50	25.79#
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	7.811	152	227380	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	1081046	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.463	164	750736	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.228	188	1718613	20.000	ng/ul	0.01
79) Chrysene-d12	21.322	240	1983835	20.000	ng/ul	0.01
88) Perylene-d12	23.721	264	1971787	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.246	96	34217	6.059	ng/uL	0.00
4) Pyridine-d5	3.681	84	69859	4.306	ng/ul	0.01
7) Phenol-d5	6.981	99	631980	30.988	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.146	67	375186	30.817	ng/ul	0.00
11) 2-Chlorophenol-d4	7.340	132	490278	32.920	ng/ul	0.00
15) 4-Methylphenol-d8	8.528	113	514664	31.907	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	256646	36.610	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143	282678	38.849	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.240	165	506545	32.123	ng/ul	0.01
31) 4-Chloroaniline-d4	10.793	131	4892m	0.208	ng/ul	0.04
46) Dimethylphthalate-d6	13.881	166	1731960	33.701	ng/ul	0.01
49) Acenaphthylene-d8	14.157	160	1978782	33.693	ng/ul	0.01
54) 4-Nitrophenol-d4	14.681	143	323347	33.244	ng/ul	0.00
60) Fluorene-d10	15.463	176	1557740	35.403	ng/ul	0.01
65) 4,6-Dinitro-2-methylph...	15.598	200	240307	26.856	ng/ul	0.01
73) Anthracene-d10	17.328	188	2591651	34.998	ng/ul	0.01
81) Pyrene-d10	19.569	212	3177209	31.899	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.569	264	3326083	34.523	ng/ul	0.02

Target Compounds Qvalue

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

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