

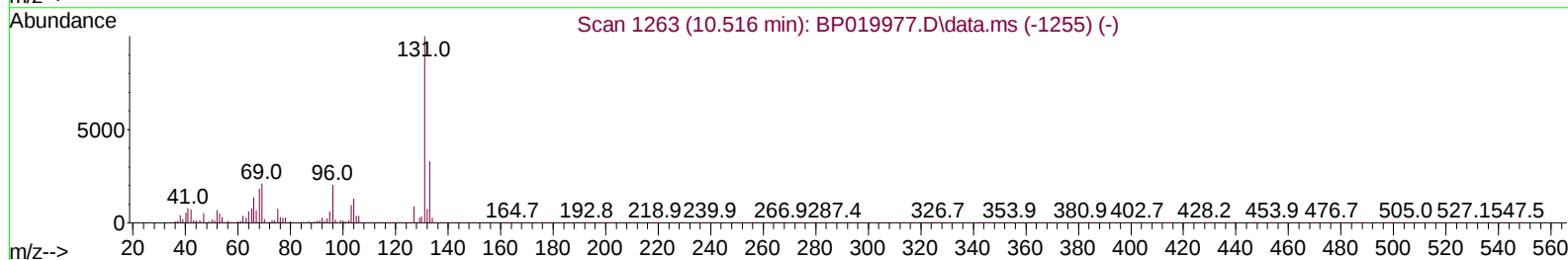
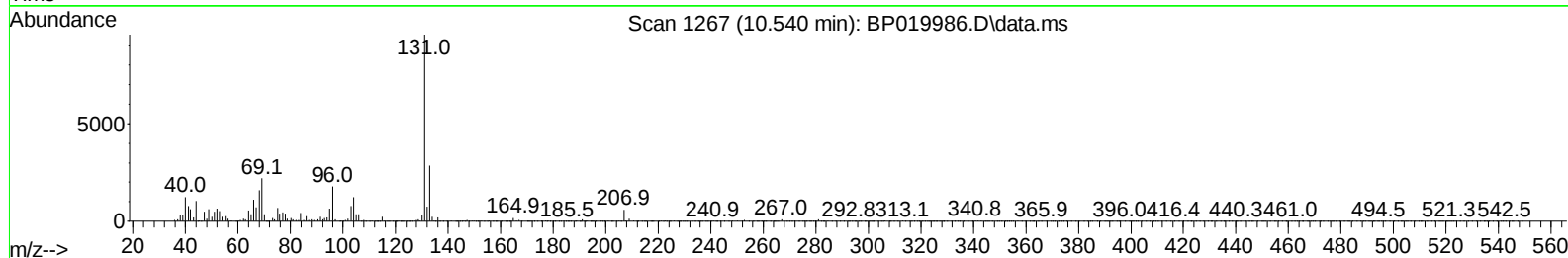
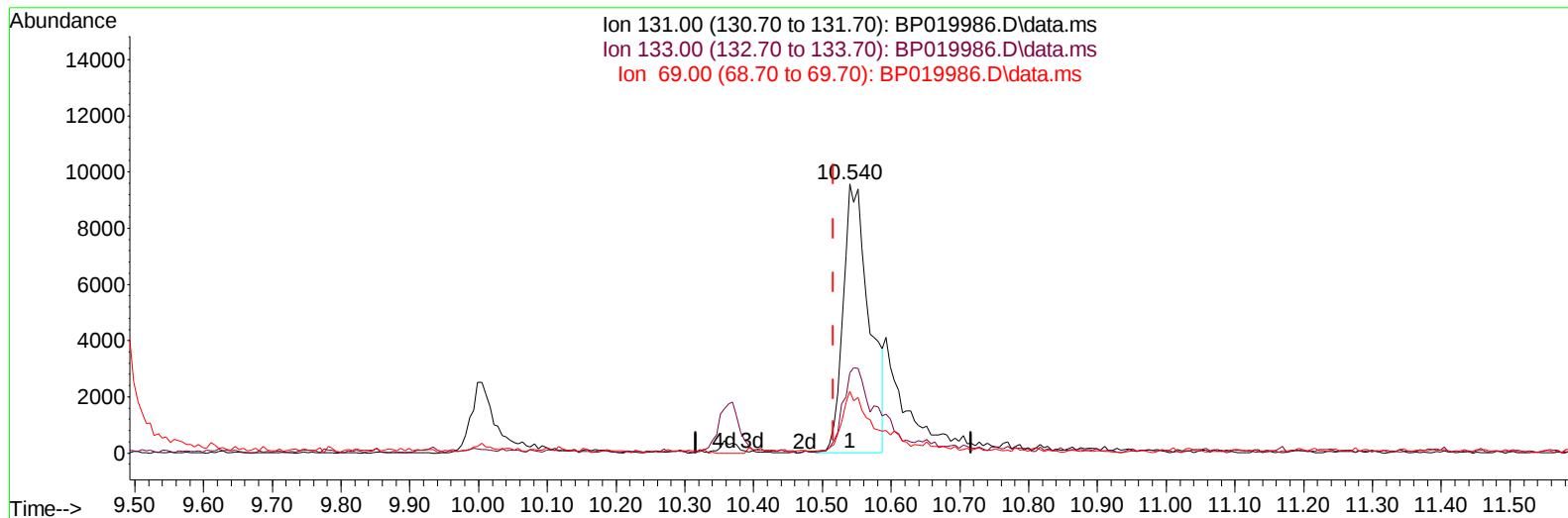
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP032524\
 Data File : BP019986.D
 Acq On : 25 Mar 2024 18:09
 Operator : MA/JU
 Sample : PB159760BL
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK760

Manual IntegrationsAPPROVED

Quant Time: Mar 25 19:06:58 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Mar 25 10:46:59 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/26/2024
 Supervised By :mohammad ahmed 03/27/2024



TIC: BP019986.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.540min (+ 0.024) 2.81 ng/u1

response 25176

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	29.88
69.00	22.40	23.00
0.00	0.00	0.00

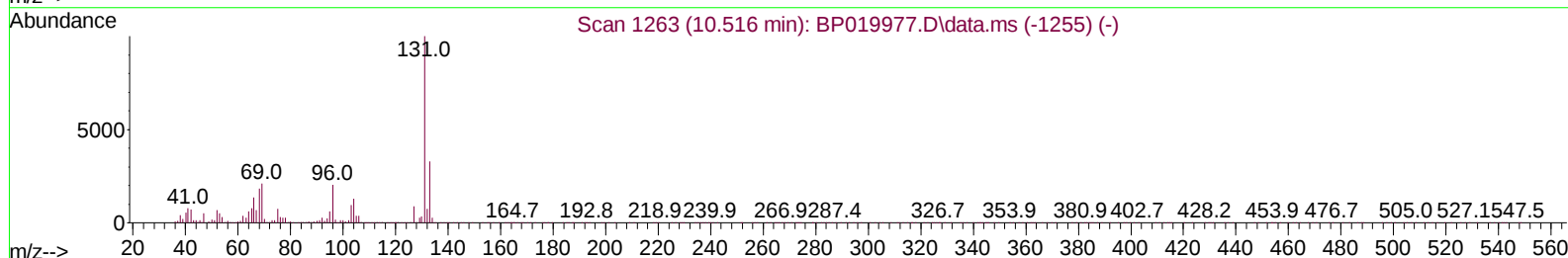
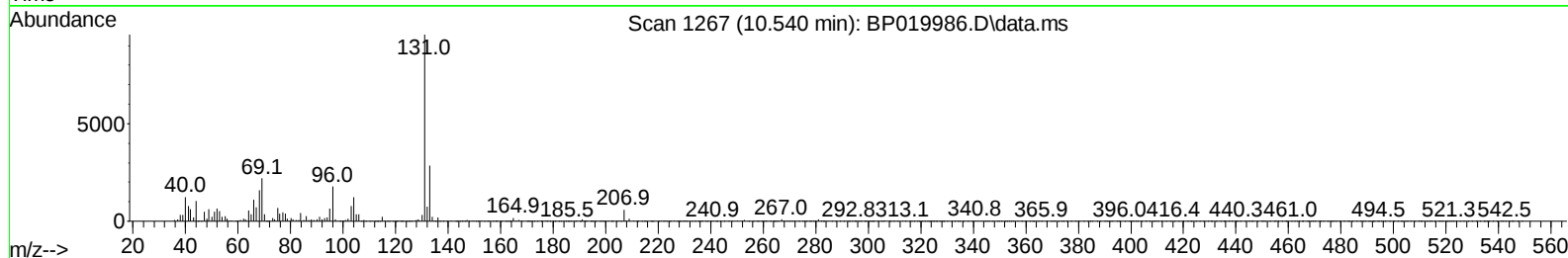
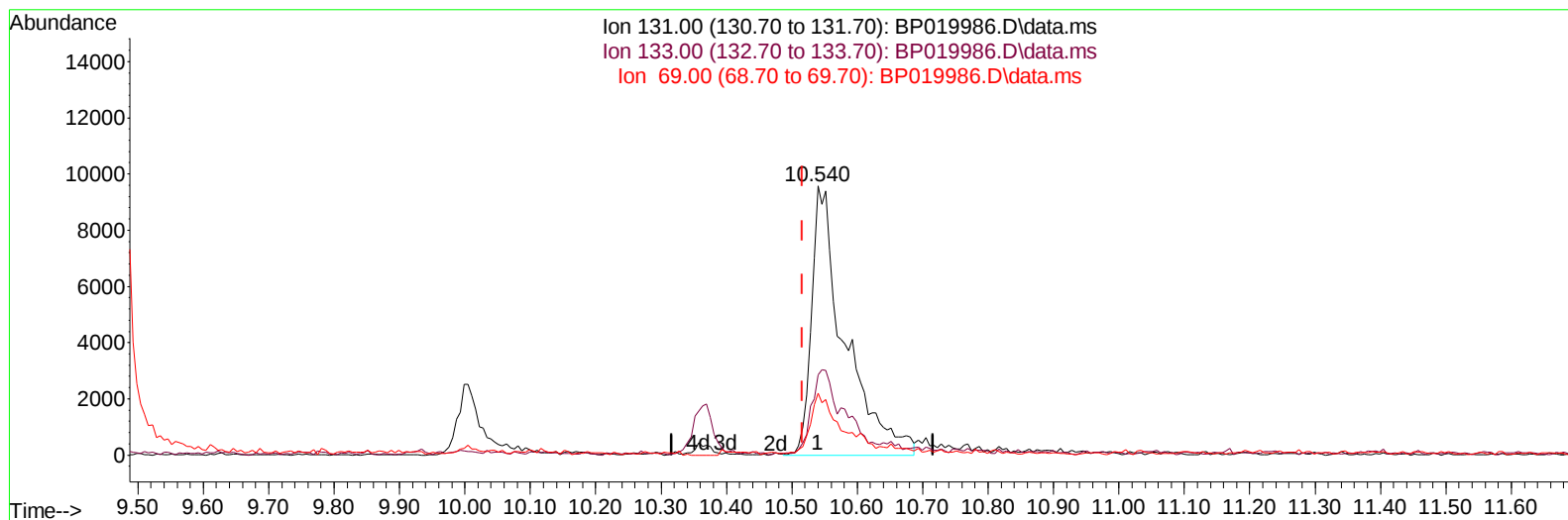
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TIC: BP019986.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.540min (+ 0.024) 3.78 ng/ul m

response 33810

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.80	29.88
69.00	22.40	23.00
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/26/2024
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Quant Time: Mar 25 19:07:28 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	111322	20.000	ng/ul	0.00
20) Naphthalene-d8	10.363	136	438148	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.234	164	292099	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.069	188	647862	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	667351	20.000	ng/ul	0.00
88) Perylene-d12	24.821	264	755415	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.240	96	16155	6.358	ng/uL	0.00
4) Pyridine-d5	3.646	84	70621	9.205	ng/ul	0.00
7) Phenol-d5	6.805	99	273855	29.168	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.964	67	184456	30.155	ng/ul	0.00
11) 2-Chlorophenol-d4	7.158	132	197573	29.302	ng/ul	0.00
15) 4-Methylphenol-d8	8.316	113	210411	28.967	ng/ul	0.01
21) Nitrobenzene-d5	8.763	128	99155	29.525	ng/ul	0.00
24) 2-Nitrophenol-d4	9.469	143	112106	29.835	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.999	165	172449	25.130	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	33810m	3.779	ng/ul	0.02
46) Dimethylphthalate-d6	13.663	166	671057	32.767	ng/ul	0.01
49) Acenaphthylene-d8	13.922	160	706690	29.539	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.493	143	92764	28.812	ng/ul	0.01
60) Fluorene-d10	15.263	176	551320	31.602	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	102793	26.507	ng/ul	-0.02
73) Anthracene-d10	17.175	188	895149	31.213	ng/ul	0.00
81) Pyrene-d10	19.563	212	1127026	28.784	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.592	264	1164330	32.083	ng/ul	-0.02

Target Compounds Qvalue

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

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