

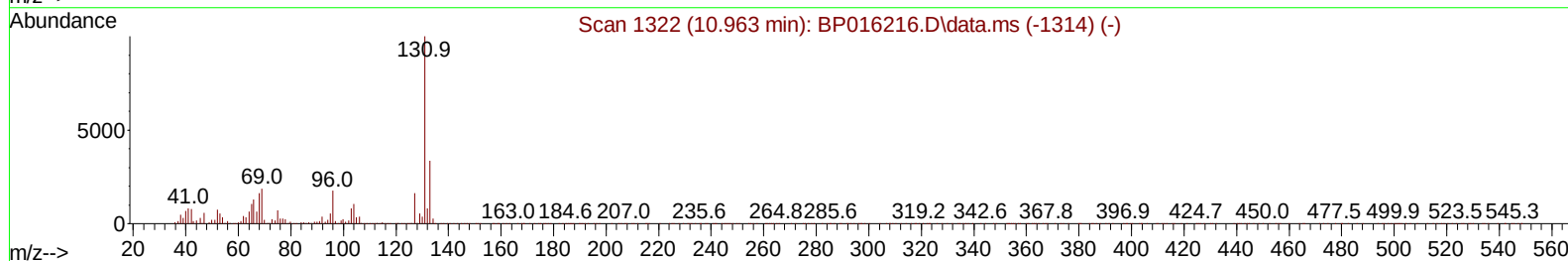
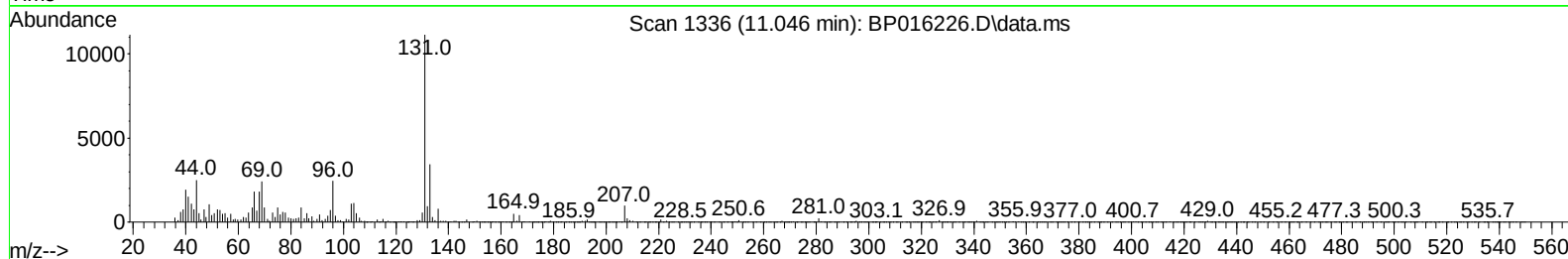
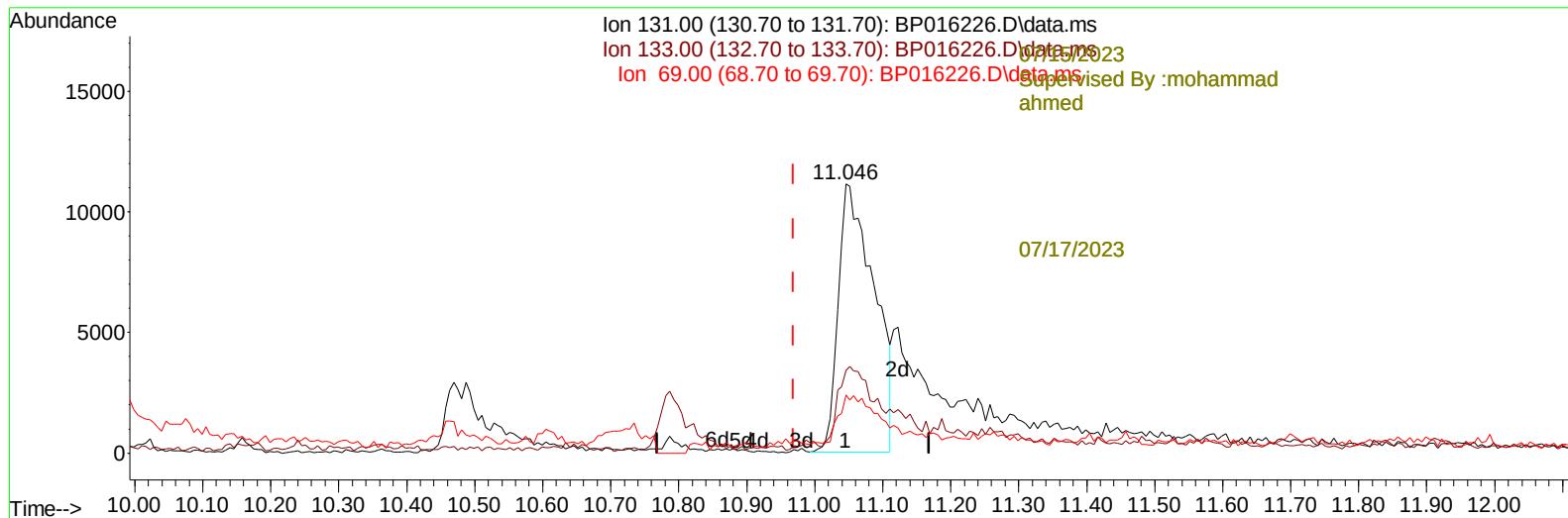
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP071423\
Data File : BP016226.D
Acq On : 14 Jul 2023 15:06
Operator : MA/JU
Sample : 03437-14
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A46G7

Manual IntegrationsAPPROVED

Quant Time: Jul 14 22:22:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jul 13 18:57:06 2023
Response via : Initial Calibration

Reviewed By :Jagrut
Upadhyay



TIC: BP016226.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.046min (+ 0.077) 2.70 ng/ul

response 40617

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.50	30.88
69.00	21.00	21.59
0.00	0.00	0.00

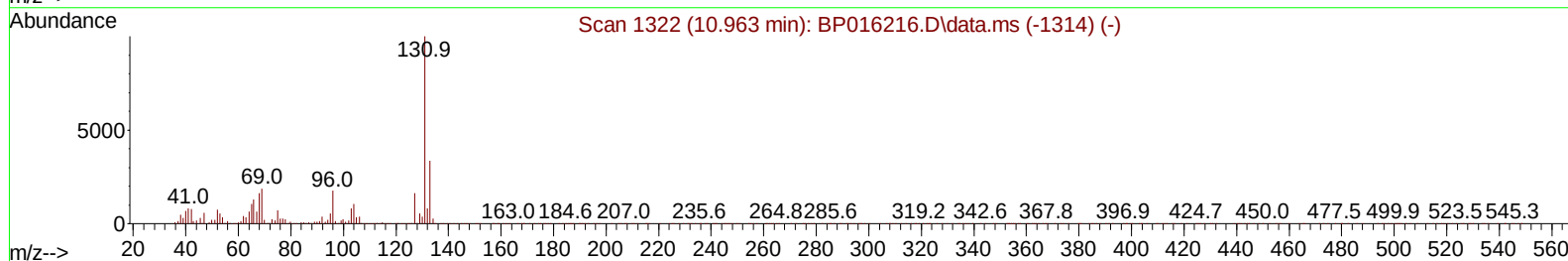
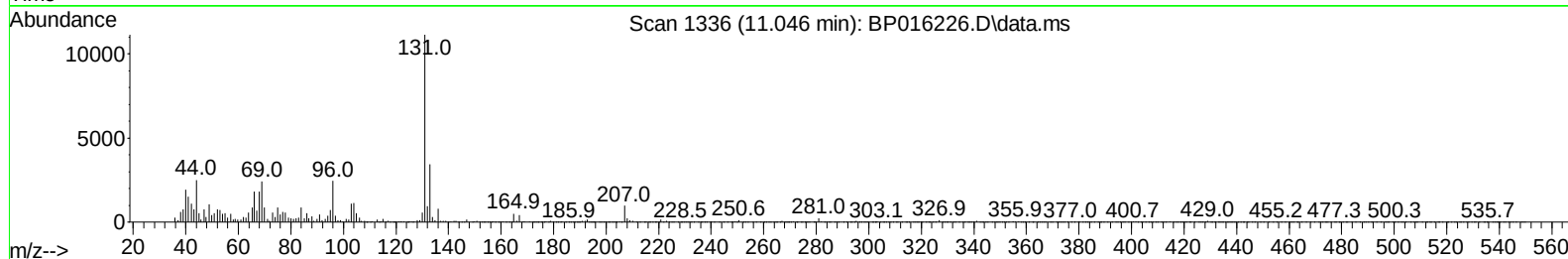
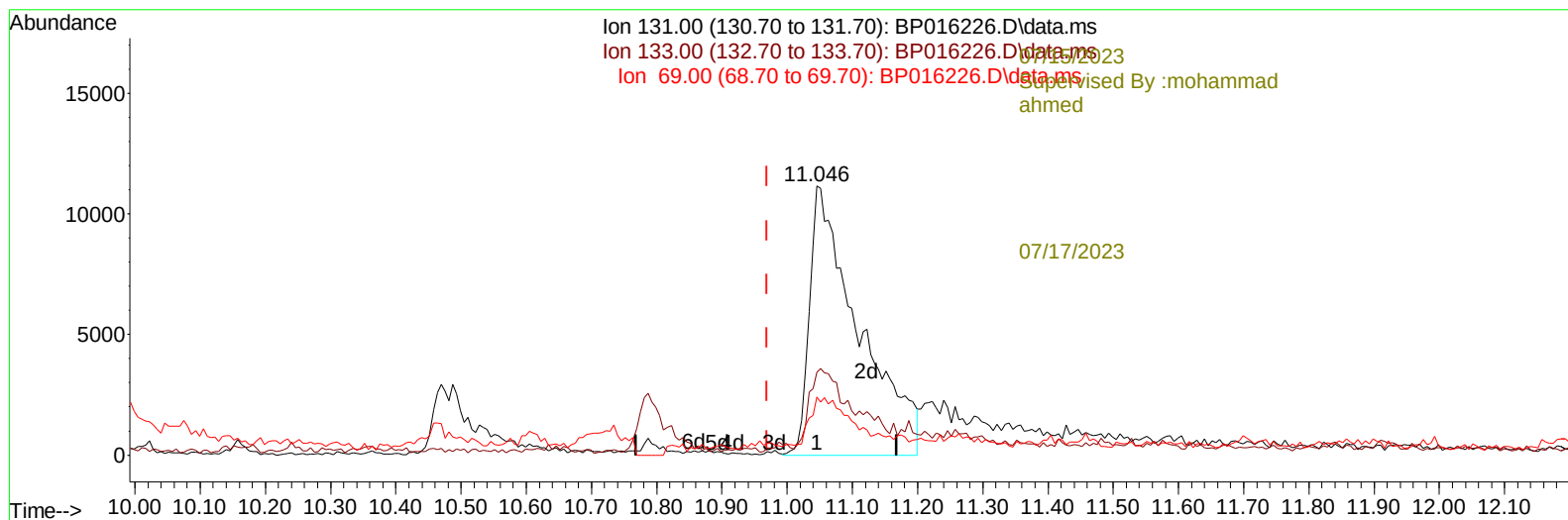
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Upadhyay



TIC: BP016226.D\data.ms

(31) 4-Chloroaniline-d4 (S)

11.046min (+ 0.077) 3.86 ng/ul m

response 57936

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	31.50	30.88
69.00	21.00	21.59
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

Quant Time: Jul 14 23:44:53 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP071223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 13 18:57:06 2023
 Response via : Initial Calibration

Reviewed By :Jagrut
 Upadhyay

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----07/15/2023							
Supervised By :mohammad							
aimed							
-----07/17/2023							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.958	152		145523	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.787	136		688100	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.616	164		451123	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.381	188		1048046	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240		1108500	20.000	ng/ul	0.00
88) Perylene-d12	23.974	264		1285824	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.281	96		28118	6.837	ng/uL	0.00
4) Pyridine-d5	3.740	84		251556	24.527	ng/ul	0.00
7) Phenol-d5	7.146	99		557043	42.379	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.287	67		360596	40.570	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.493	132		415414	43.780	ng/ul	0.00
15) 4-Methylphenol-d8	8.716	113		215527	20.355	ng/ul	0.00
21) Nitrobenzene-d5	9.158	128		206522	43.504	ng/ul	0.00
24) 2-Nitrophenol-d4	9.887	143		233616	43.178	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.469	165		390881	41.719	ng/ul	0.01
31) 4-Chloroaniline-d4	11.046	131		57936m	3.858	ng/ul	0.08
46) Dimethylphthalate-d6	14.034	166		1420179	41.270	ng/ul	0.00
49) Acenaphthylene-d8	14.310	160		1548080	41.049	ng/ul	0.00
54) 4-Nitrophenol-d4	14.934	143		174941	24.910	ng/ul	0.00
60) Fluorene-d10	15.616	176		1246814	43.293	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.798	200		167795	27.734	ng/ul	0.00
73) Anthracene-d10	17.487	188		1451598	31.664	ng/ul	0.00
81) Pyrene-d10	19.716	212		2496635	43.405	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.804	264		1832499	29.677	ng/ul	-0.01
Target Compounds							
						Qvalue	
16) Acetophenone	8.834	105		59461	3.405	ng/ul	97
86) Bis(2-ethylhexyl)phtha...	21.333	149		132763	2.870	ng/ul	97

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

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