

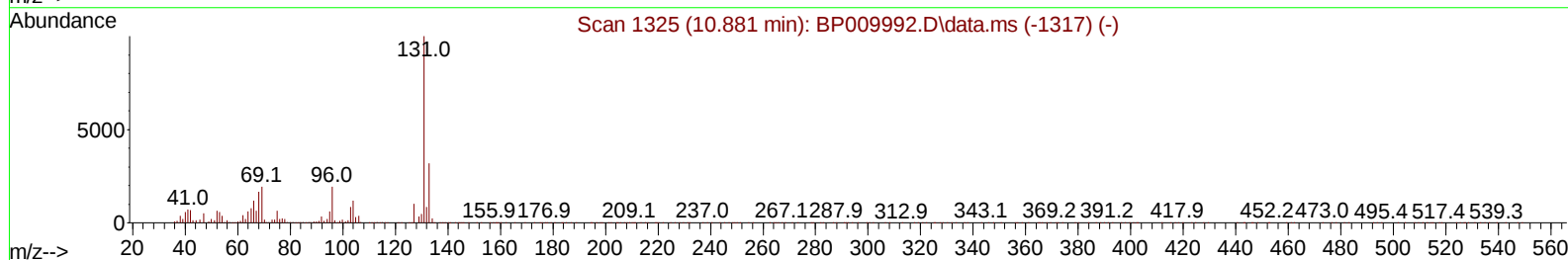
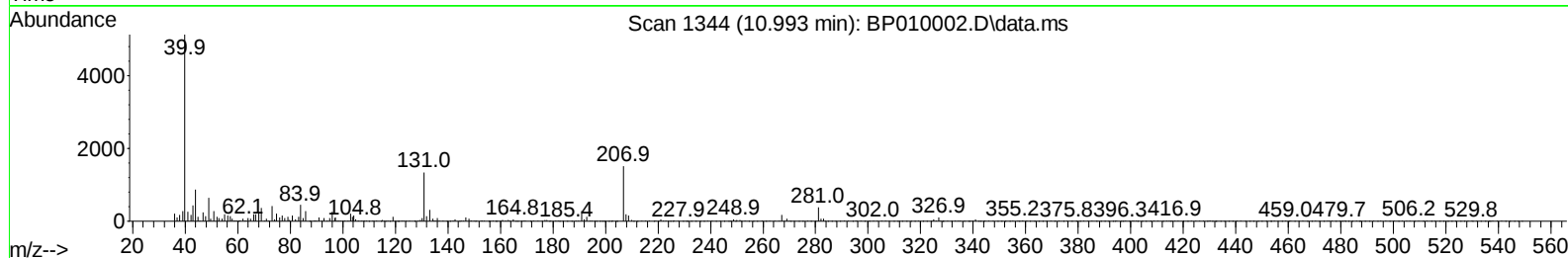
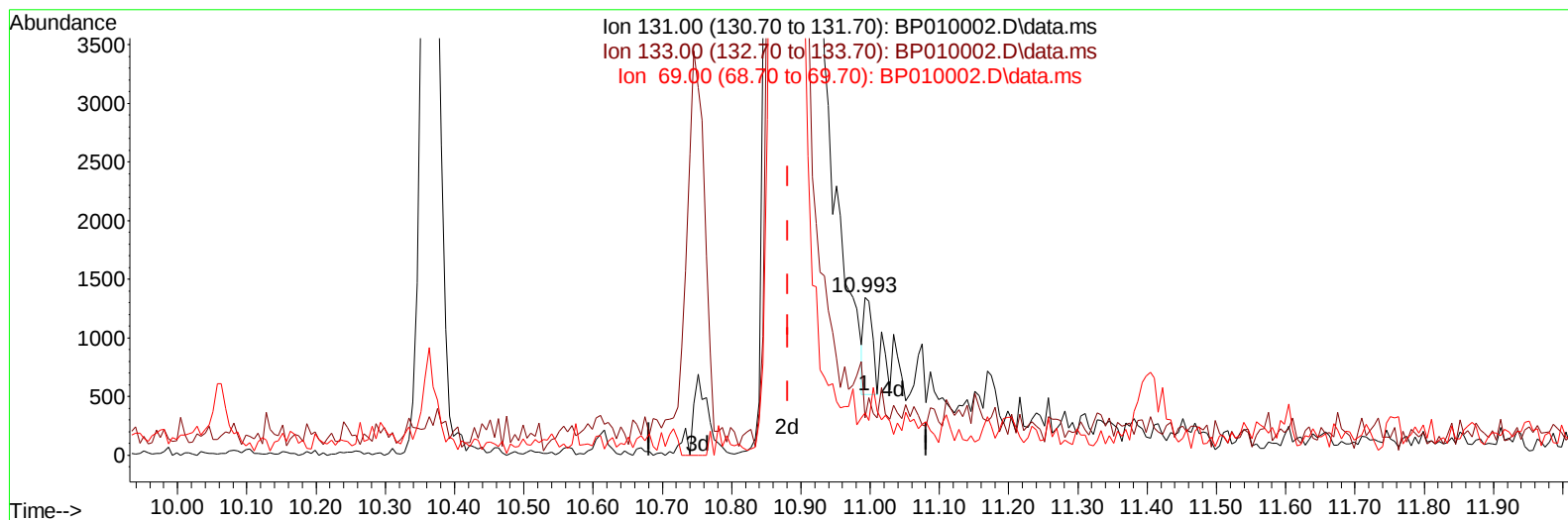
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP010002.D
 Acq On : 21 Apr 2022 06:10
 Operator : CG/JU
 Sample : N2473-02
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J24

Manual IntegrationsAPPROVED

Quant Time: Apr 21 06:54:54 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/21/2022
 Supervised By :Yogesh Patel 04/25/2022



TIC: BP010002.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.993min (+ 0.112) 0.05 ng/u1

response 739

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	23.40
69.00	26.10	26.60
0.00	0.00	0.00

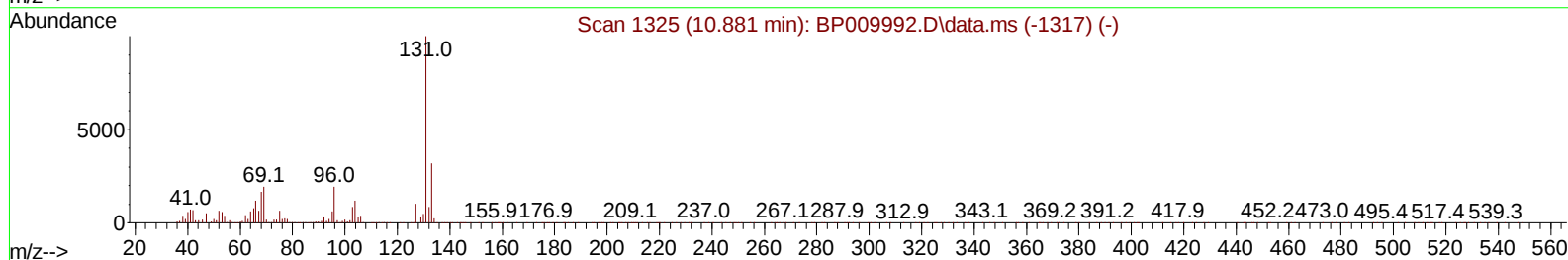
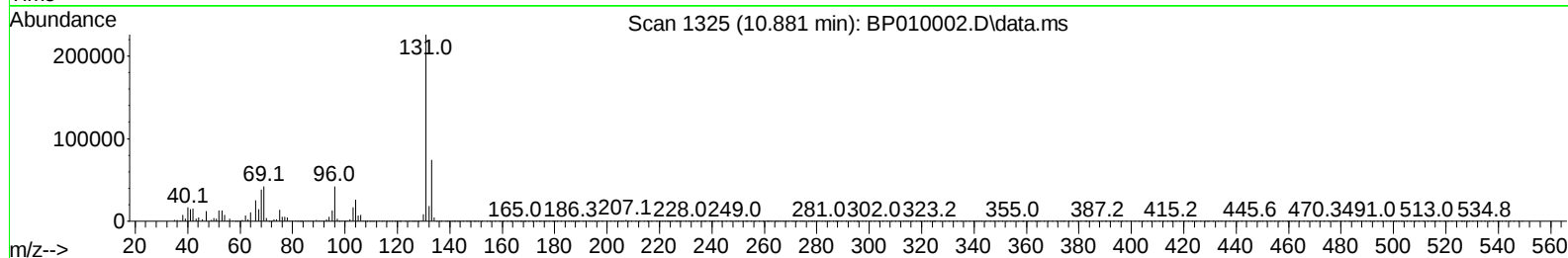
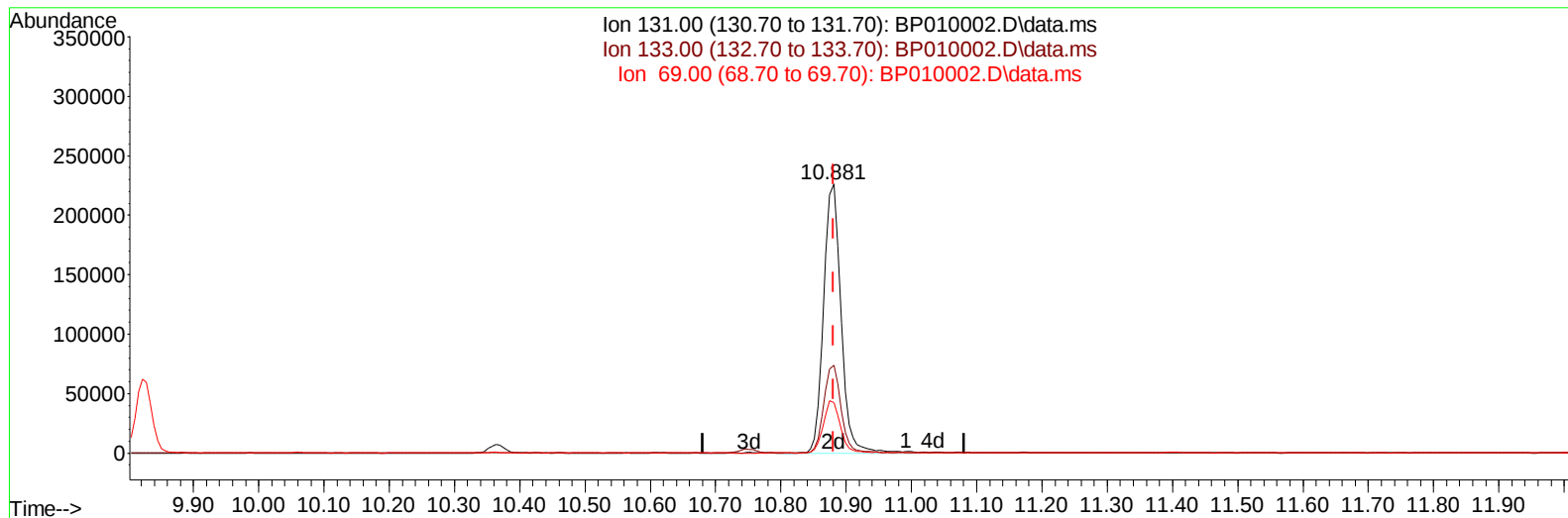
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TIC: BP010002.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.881min (-0.000) 25.81 ng/ul m

response 414145

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	32.79
69.00	26.10	18.66#
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	7.940	152	146722	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	680960	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.575	164	485867	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1076450	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	1020100	20.000	ng/ul	# 0.00
88) Perylene-d12	23.851	264	836560	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	15706	4.777	ng/uL	0.00
4) Pyridine-d5	3.793	84	86187	9.508	ng/ul	0.00
7) Phenol-d5	7.099	99	91707	7.119	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	265960	34.681	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	257028	26.453	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	190197	17.688	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	178956	36.443	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	188795	35.068	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	330314	29.320	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	414145m	25.810	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1347846	35.316	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1475856	32.641	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	44564	6.438	ng/ul	0.00
60) Fluorene-d10	15.563	176	1187421	36.984	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	211044	33.095	ng/ul	-0.01
73) Anthracene-d10	17.422	188	1986425	38.542	ng/ul	0.00
81) Pyrene-d10	19.651	212	2464179	44.230	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.692	264	1753229	40.244	ng/ul	-0.02

Target Compounds Qvalue

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

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