

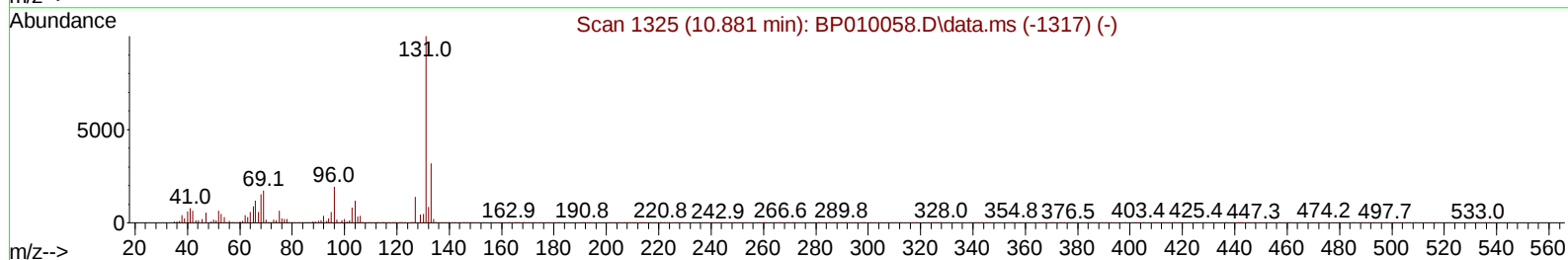
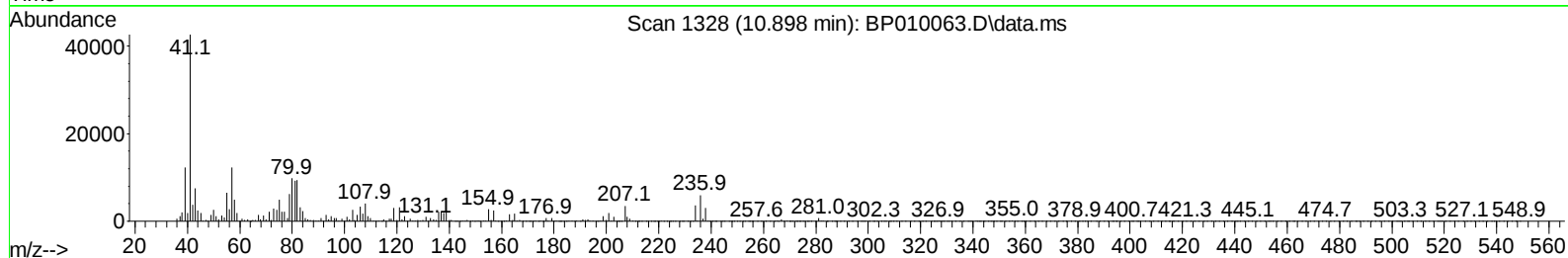
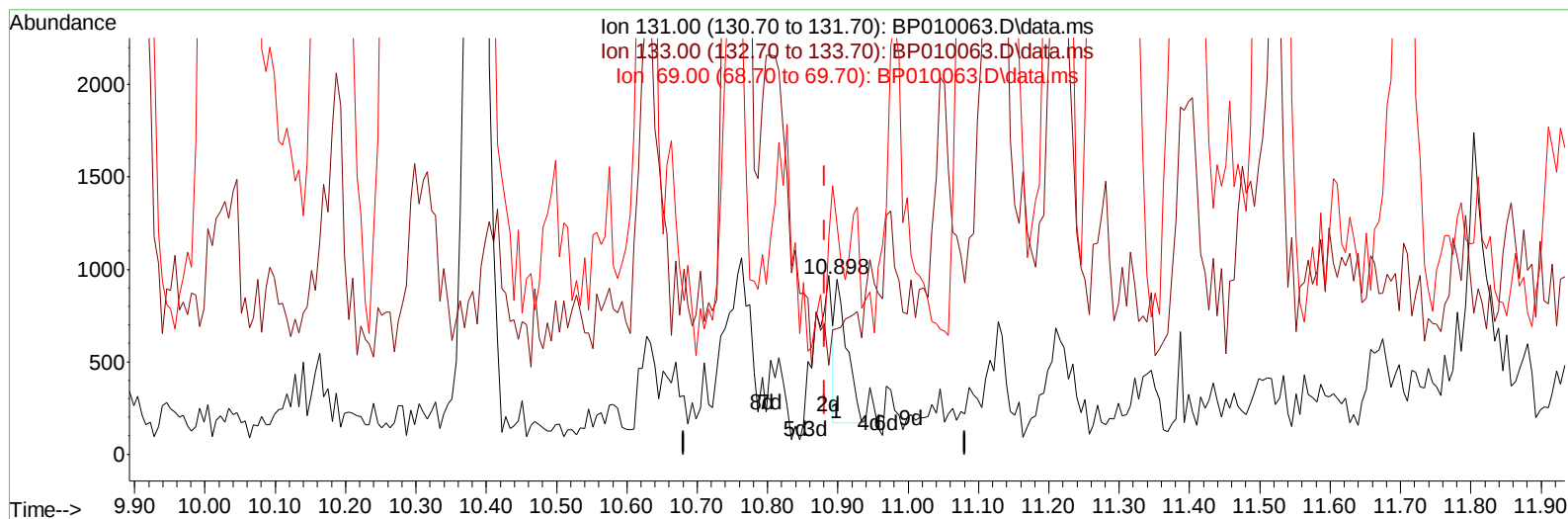
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010063.D
 Acq On : 22 Apr 2022 20:00
 Operator : CG/JU
 Sample : N2316-14RE
 Misc :
 ALS Vial : 59 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 ETNR4RE

Manual IntegrationsAPPROVED

Quant Time: Apr 22 23:30:47 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/25/2022
 Supervised By :mohammad ahmed 04/26/2022



TIC: BP010063.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.898min (+ 0.018) 0.04 ng/ul

response 904

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	71.73#
69.00	26.10	131.01#
0.00	0.00	0.00

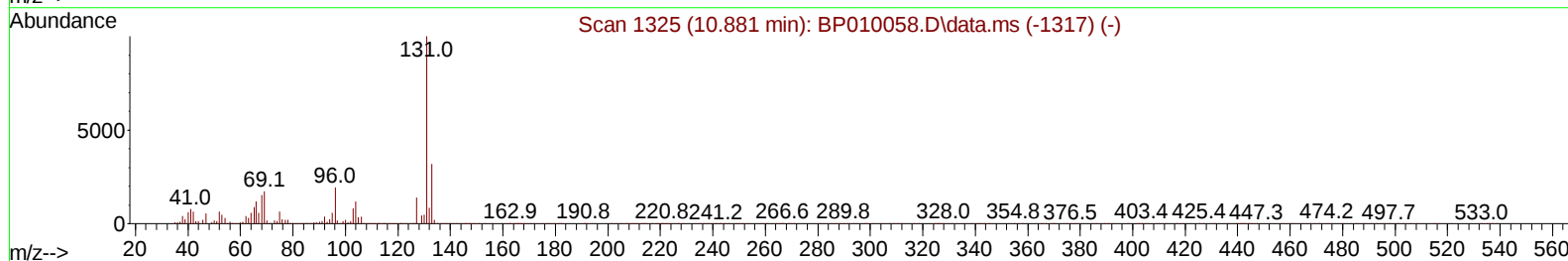
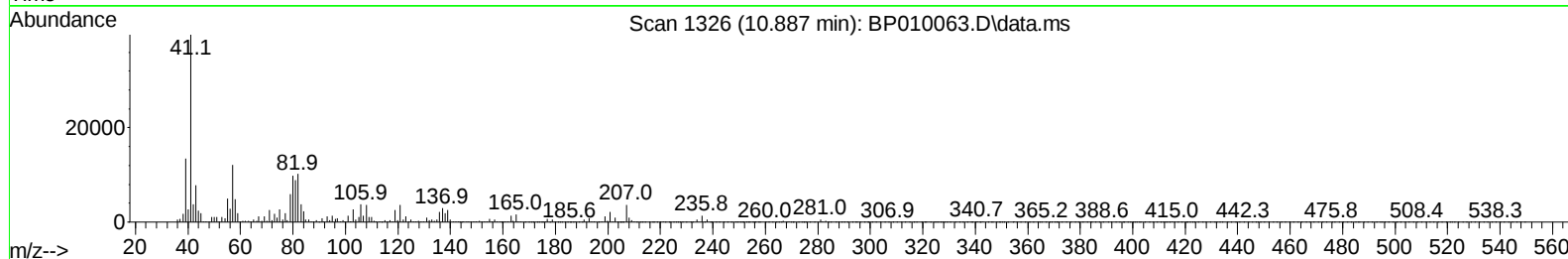
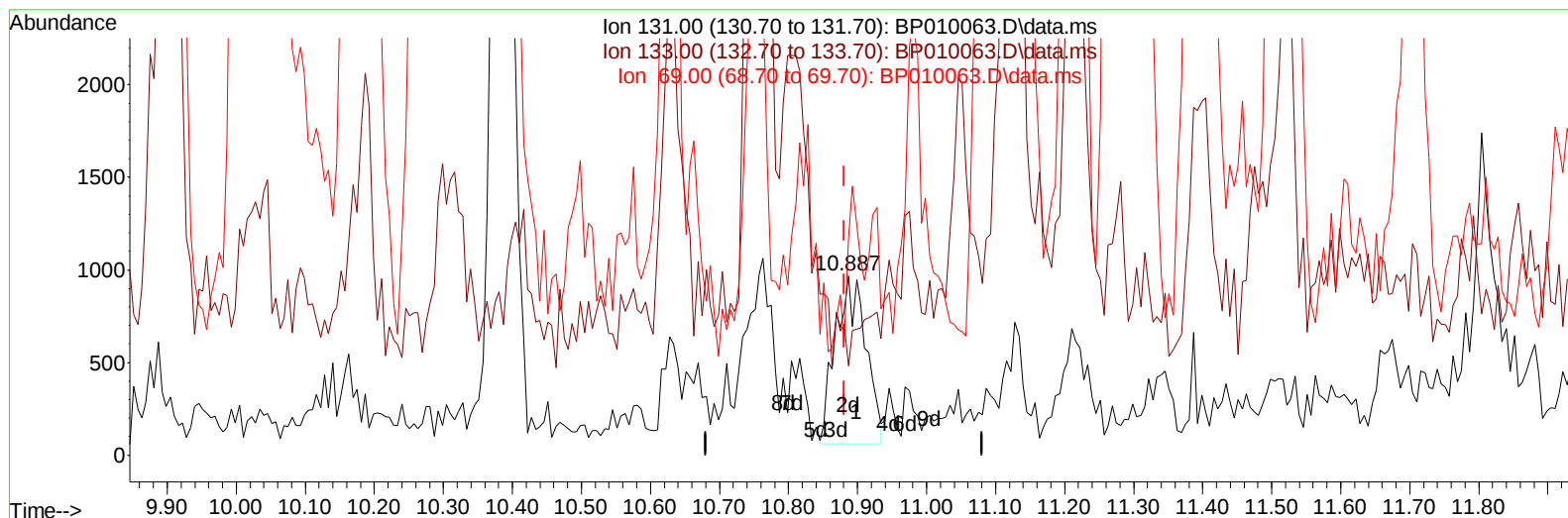
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(31) 4-Chloroaniline-d4 (S)

10.887min (+ 0.006) 0.14 ng/ul m

response 2769

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	49.64#
69.00	26.10	121.36#
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	198191	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136	842594	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.581	164	527432	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.357	188	1020718	20.000	ng/ul	# 0.03
79) Chrysene-d12	21.416	240	686179	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264	96248	20.000	ng/ul	# 0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.375	96	18870	4.152	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.110	99	419879	23.291	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.269	67	280293	25.912	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	356693	26.717	ng/ul	0.00
15) 4-Methylphenol-d8	8.657	113	339596	22.737	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	188501	28.128	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	205037	28.014	ng/ul	0.02
28) 2,4-Dichlorophenol-d3	10.387	165	504828	35.431	ng/ul	0.02
31) 4-Chloroaniline-d4	10.887	131	2769m	0.136	ng/ul	0.00
46) Dimethylphthalate-d6	13.992	166	1105356	26.941	ng/ul	0.01
49) Acenaphthylene-d8	14.281	160	1266118	25.498	ng/ul	0.02
54) 4-Nitrophenol-d4	14.775	143	119987	15.886	ng/ul	0.01
60) Fluorene-d10	15.569	176	962173	27.263	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	112892	17.491	ng/ul	0.00
73) Anthracene-d10	17.457	188	1255929	25.433	ng/ul	0.04
81) Pyrene-d10	19.669	212	1265879	32.439	ng/ul	0.01
92) Benzo(a)pyrene-d12	23.715	264	206903	41.004	ng/ul	0.02
Target Compounds						
					Qvalue	
43) 1,1'-Biphenyl	13.410	154	44013	1.131	ng/ul#	82
90) Benzo(b)fluoranthene	23.133	252	11402	1.762	ng/ul#	5

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

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