

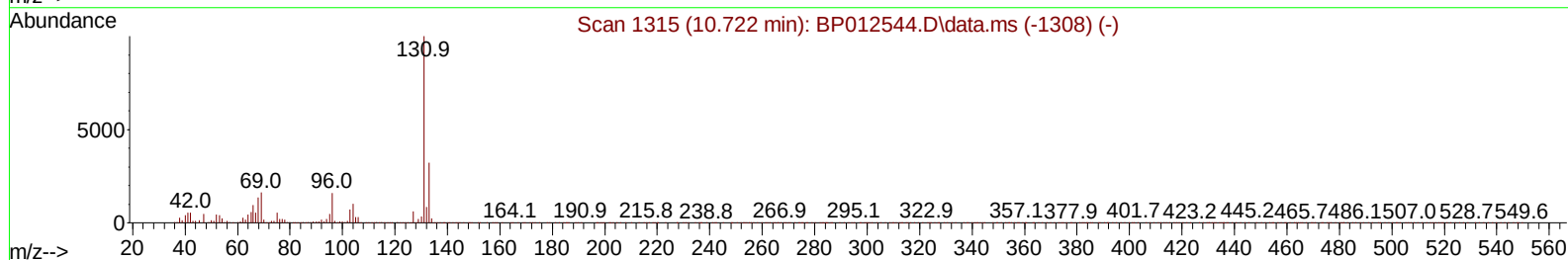
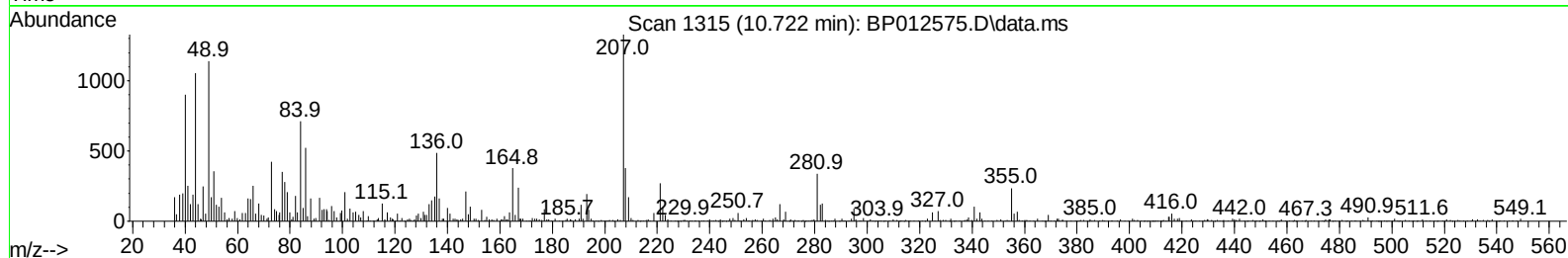
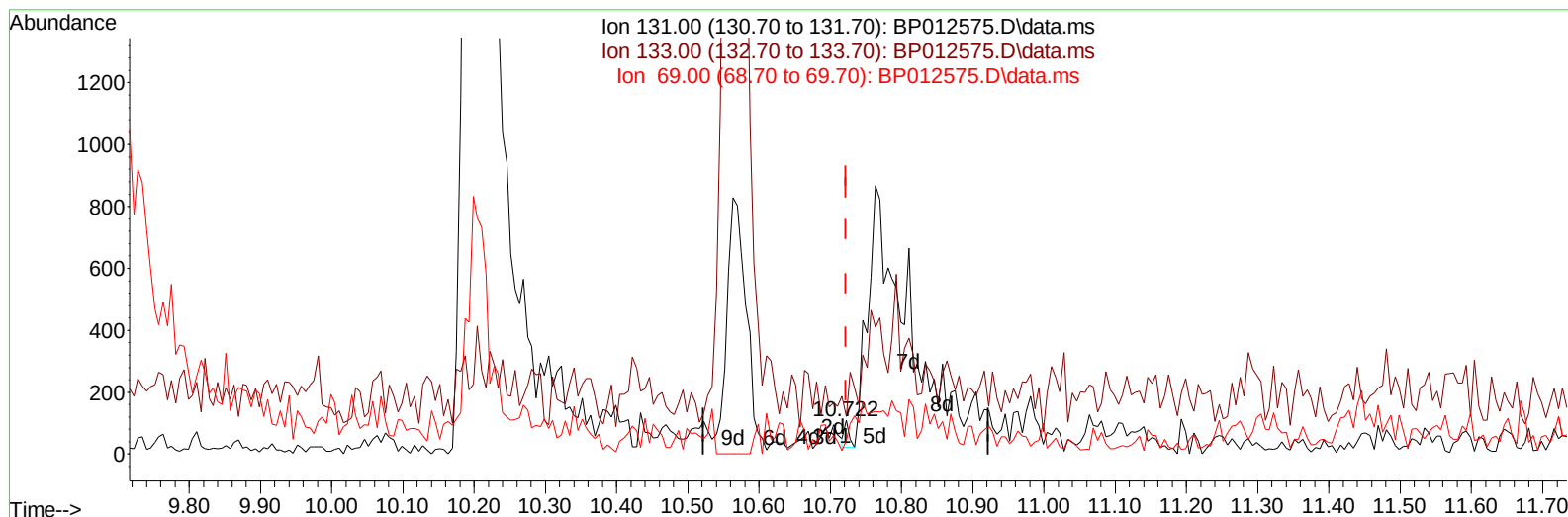
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110822\
 Data File : BP012575.D
 Acq On : 09 Nov 2022 05:29
 Operator : CG/JU
 Sample : PB148838BL
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK838

Manual IntegrationsAPPROVED

Quant Time: Nov 09 06:19:26 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP102522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 08 22:40:23 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/09/2022
 Supervised By :mohammad ahmed 11/14/2022



TIC: BP012575.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.722min (-0.000) 0.00 ng/ul

response 40

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	111.01#
69.00	16.50	39.45#
0.00	0.00	0.00

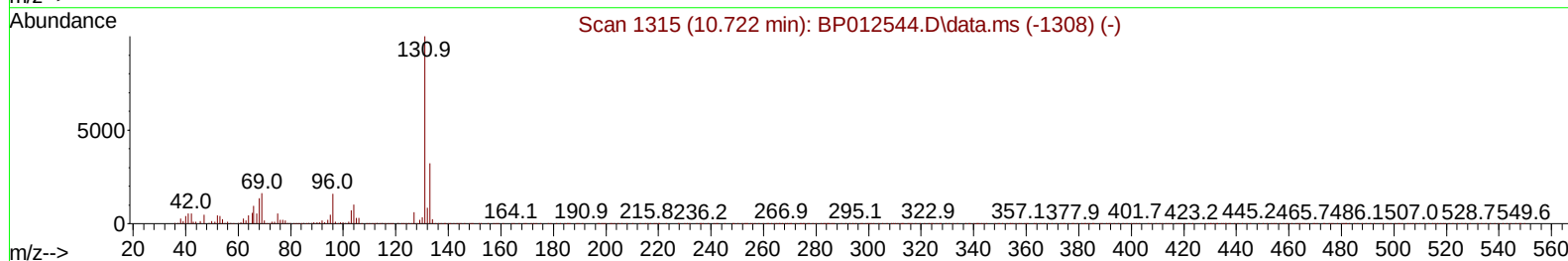
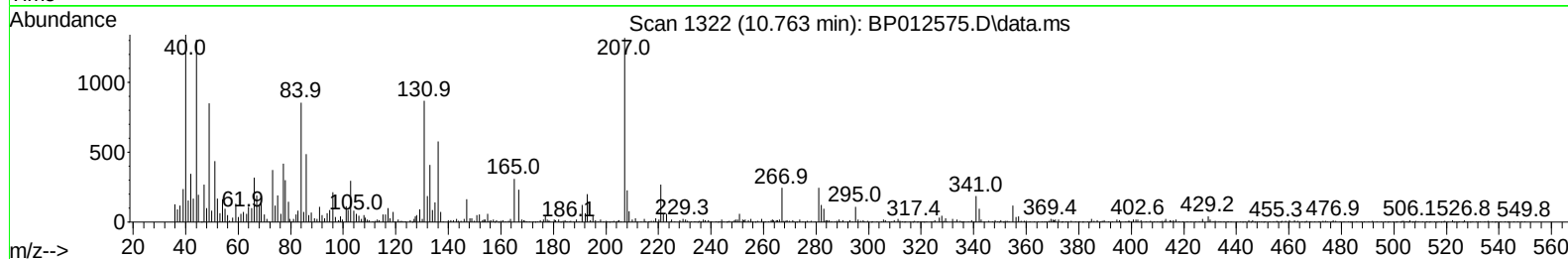
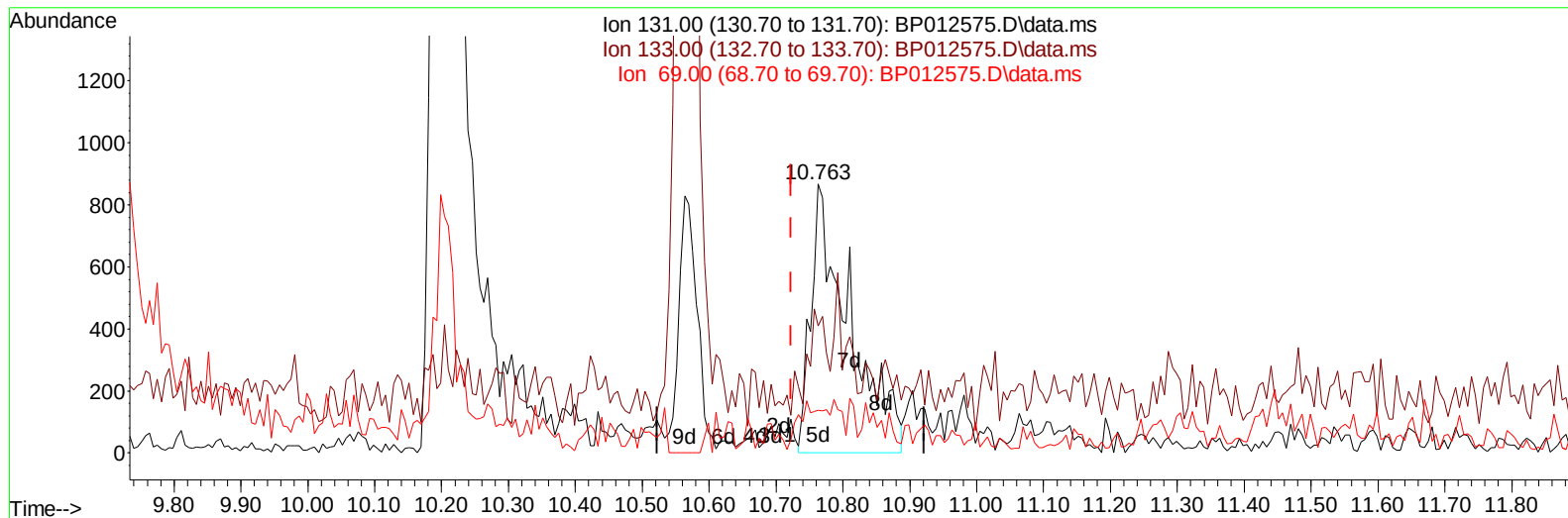
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(31) 4-Chloroaniline-d4 (S)

10.763min (+ 0.041) 0.15 ng/ul m

response 3428

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	47.17#
69.00	16.50	15.80
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	7.763	152	210070	20.000	ng/ul	0.00
20) Naphthalene-d8	10.563	136	1042413	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	766186	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	1748835	20.000	ng/ul	0.00
79) Chrysene-d12	21.280	240	1619723	20.000	ng/ul	0.00
88) Perylene-d12	23.657	264	1463609	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	32567	6.242	ng/uL	0.00
4) Pyridine-d5	3.664	84	77271	5.156	ng/ul	0.01
7) Phenol-d5	6.958	99	675912	35.873	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.105	67	382015	33.963	ng/ul	0.00
11) 2-Chlorophenol-d4	7.305	132	514260	37.375	ng/ul	0.00
15) 4-Methylphenol-d8	8.493	113	558590	37.484	ng/ul	0.00
21) Nitrobenzene-d5	8.940	128	273299	40.430	ng/ul	0.00
24) 2-Nitrophenol-d4	9.657	143	311122	44.343	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	568325	37.376	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	3428m	0.151	ng/ul	0.04
46) Dimethylphthalate-d6	13.840	166	1976765	37.689	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	2148223	35.840	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	311490	31.379	ng/ul	0.00
60) Fluorene-d10	15.416	176	1741706	38.786	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	311458	34.207	ng/ul	0.00
73) Anthracene-d10	17.281	188	2887746	38.323	ng/ul	0.00
81) Pyrene-d10	19.527	212	3322381	40.855	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264	2722643	38.071	ng/ul	0.00

Target Compounds Qvalue

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

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