

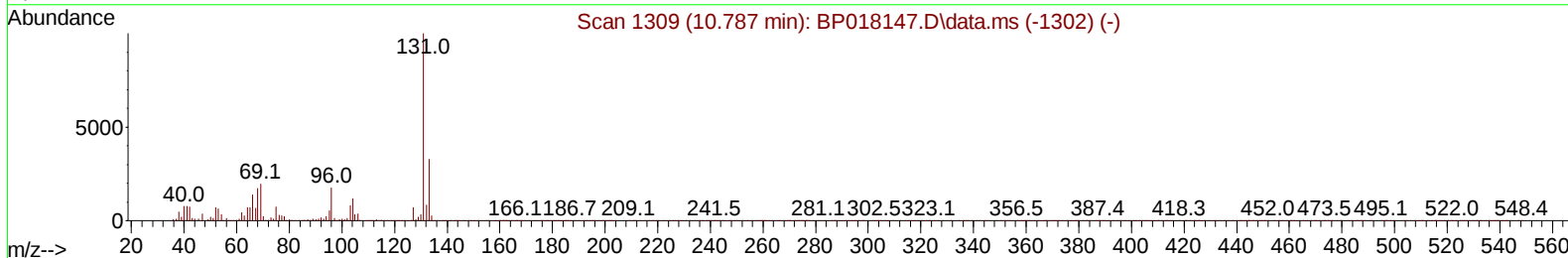
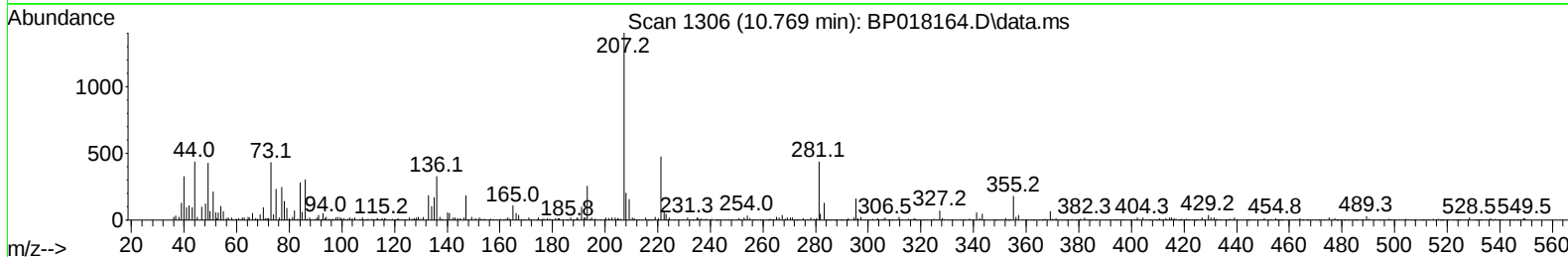
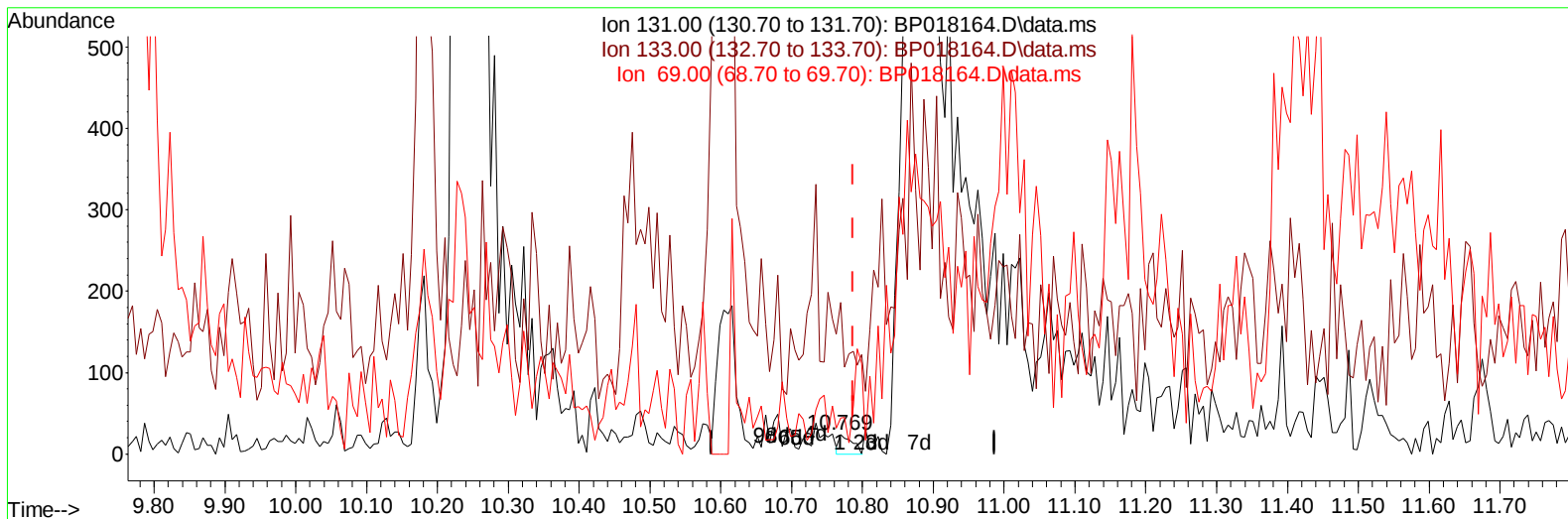
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111423\
 Data File : BP018164.D
 Acq On : 14 Nov 2023 20:06
 Operator : MA/JU
 Sample : 05337-10
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDJ5

Manual IntegrationsAPPROVED

Quant Time: Nov 15 00:40:20 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/16/2023
 Supervised By :mohammad ahmed 11/16/2023



TIC: BP018164.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.769min (-0.018) 0.01 ng/u1

response 31

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	775.00#
69.00	20.00	183.33#
0.00	0.00	0.00

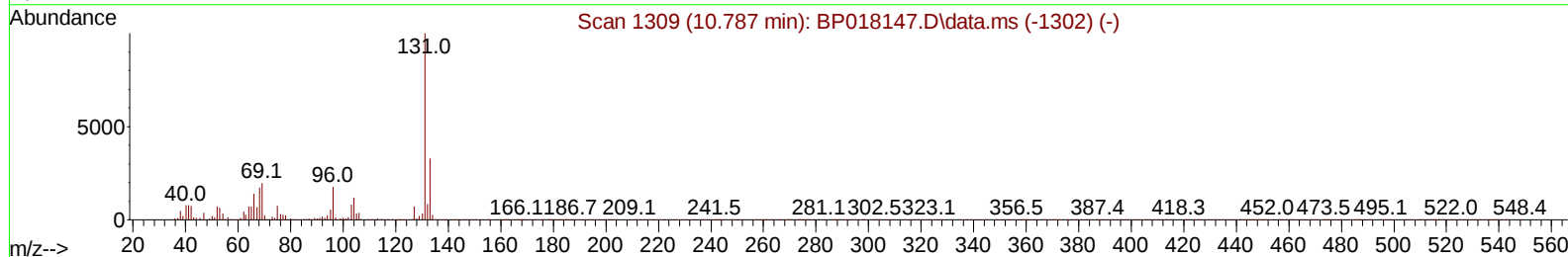
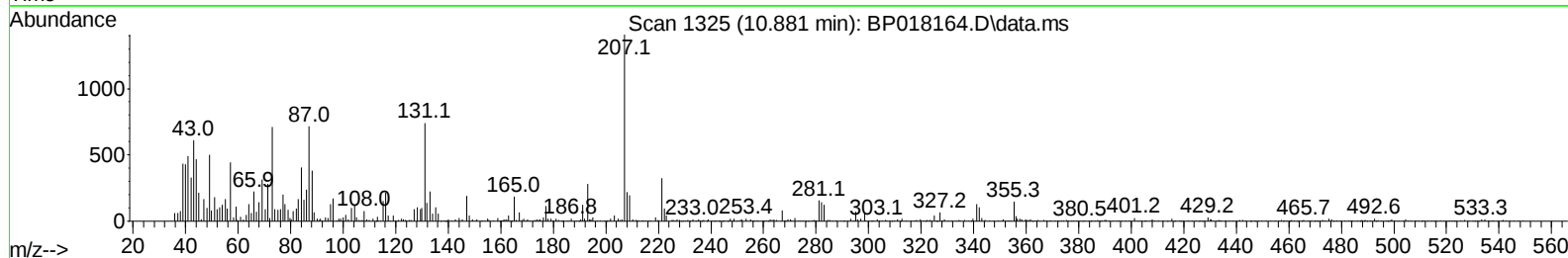
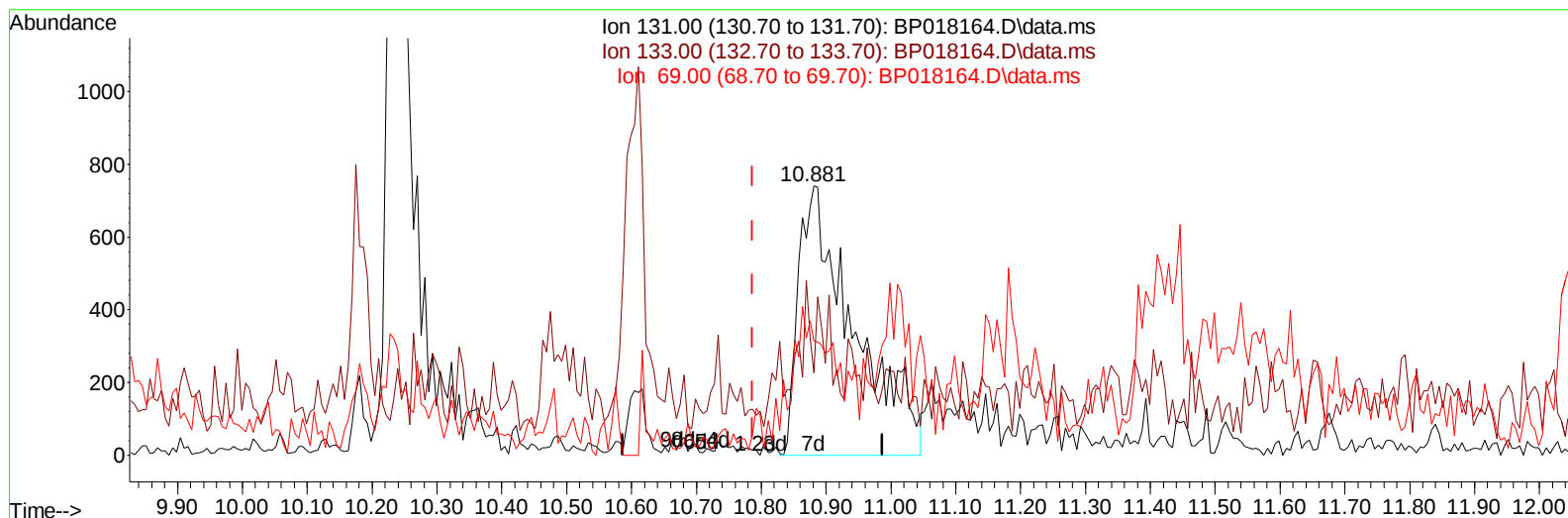
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(31) 4-Chloroaniline-d4 (S)

10.881min (+ 0.094) 0.87 ng/ul m

response 4401

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	32.20	30.54
69.00	20.00	42.43#
0.00	0.00	0.00

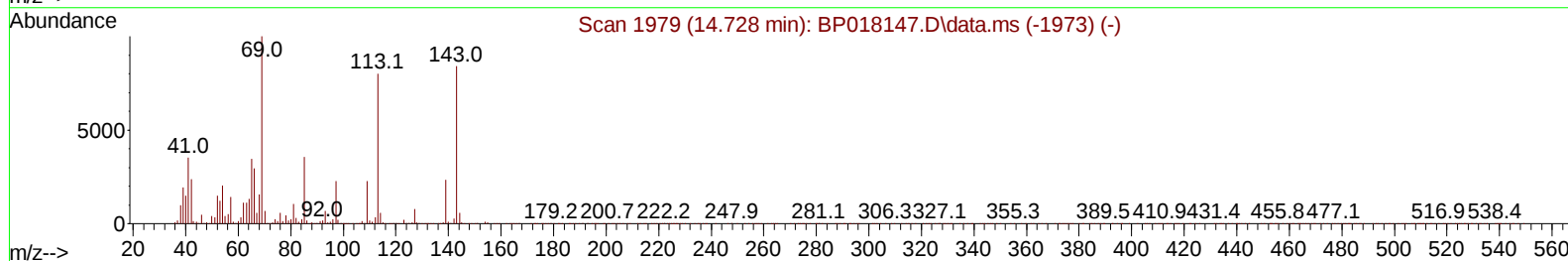
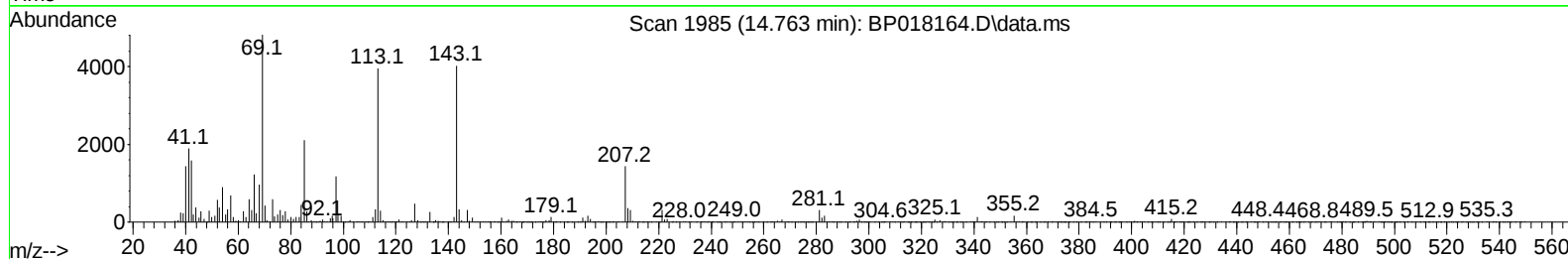
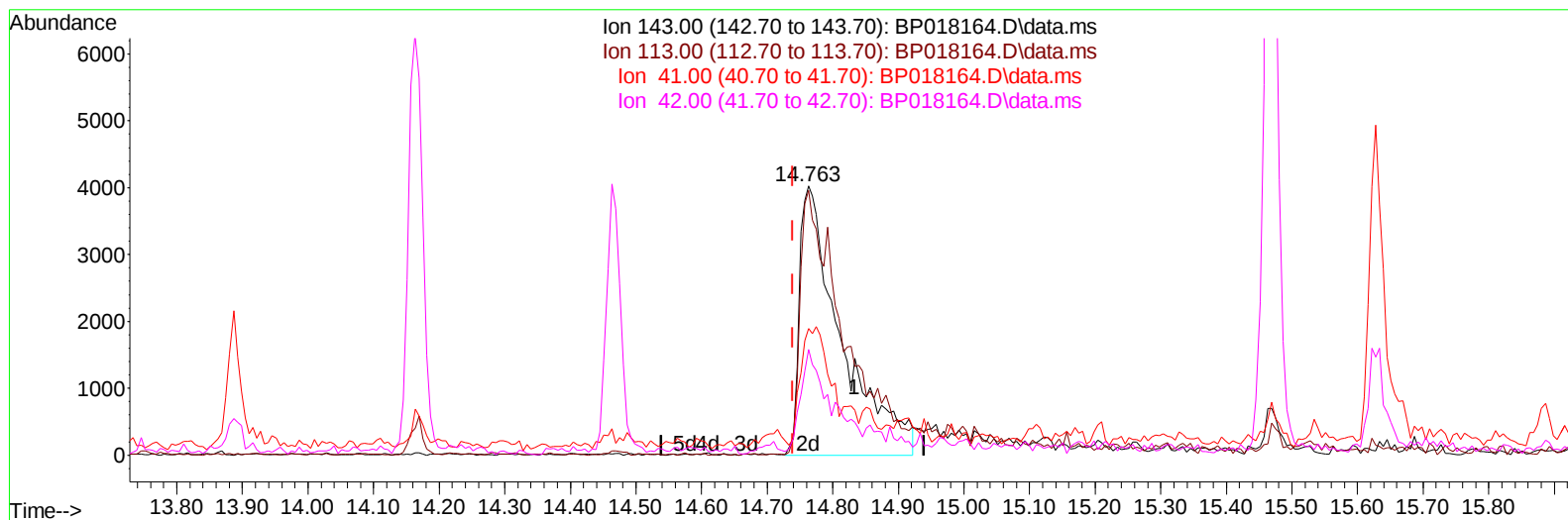
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TIC: BP018164.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.024) 10.44 ng/ul m

response 17828

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.00	98.43
41.00	47.70	47.08
42.00	30.10	39.25#

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Instrument :
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Manual IntegrationsAPPROVED

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.787	152	50281	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	201927	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.469	164	123881	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	279508	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	287658	20.000	ng/ul	-0.01
88) Perylene-d12	23.851	264	317825	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	7208	5.057	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.958	99	61639	12.698	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	125170	43.877	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	145383	37.820	ng/ul	0.00
15) 4-Methylphenol-d8	8.510	113	109161	27.443	ng/ul	0.00
21) Nitrobenzene-d5	8.993	128	74383	40.887	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	96225	46.688	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.234	165	158319	43.099	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	4401m	0.874	ng/ul	0.09
46) Dimethylphthalate-d6	13.887	166	548575	47.993	ng/ul	-0.01
49) Acenaphthylene-d8	14.163	160	513882	41.277	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	17828m	10.438	ng/ul	0.02
60) Fluorene-d10	15.469	176	439470	46.569	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	73825	36.536	ng/ul	-0.04
73) Anthracene-d10	17.345	188	749260	49.564	ng/ul	-0.01
81) Pyrene-d10	19.604	212	973938	50.688	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.686	264	1010582	54.683	ng/ul	-0.02

Target Compounds Qvalue

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

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