

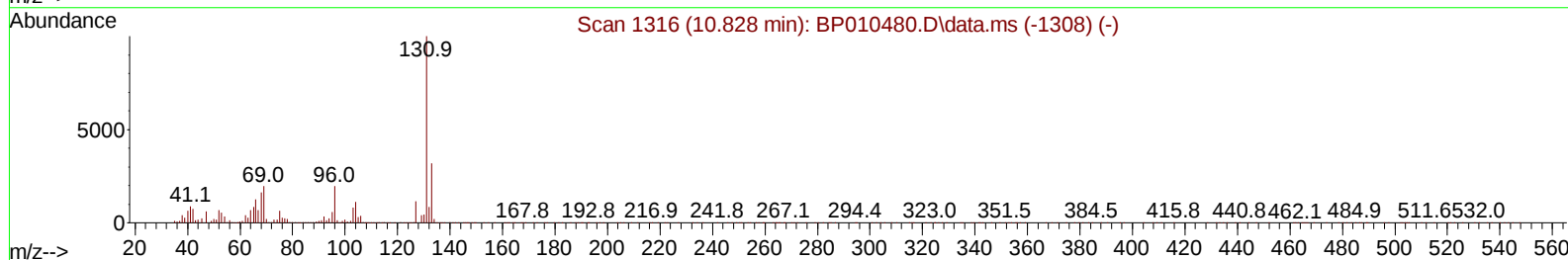
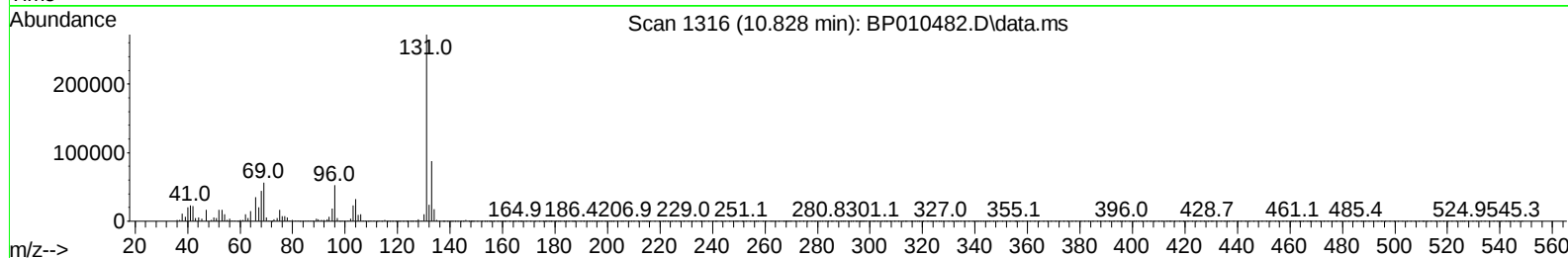
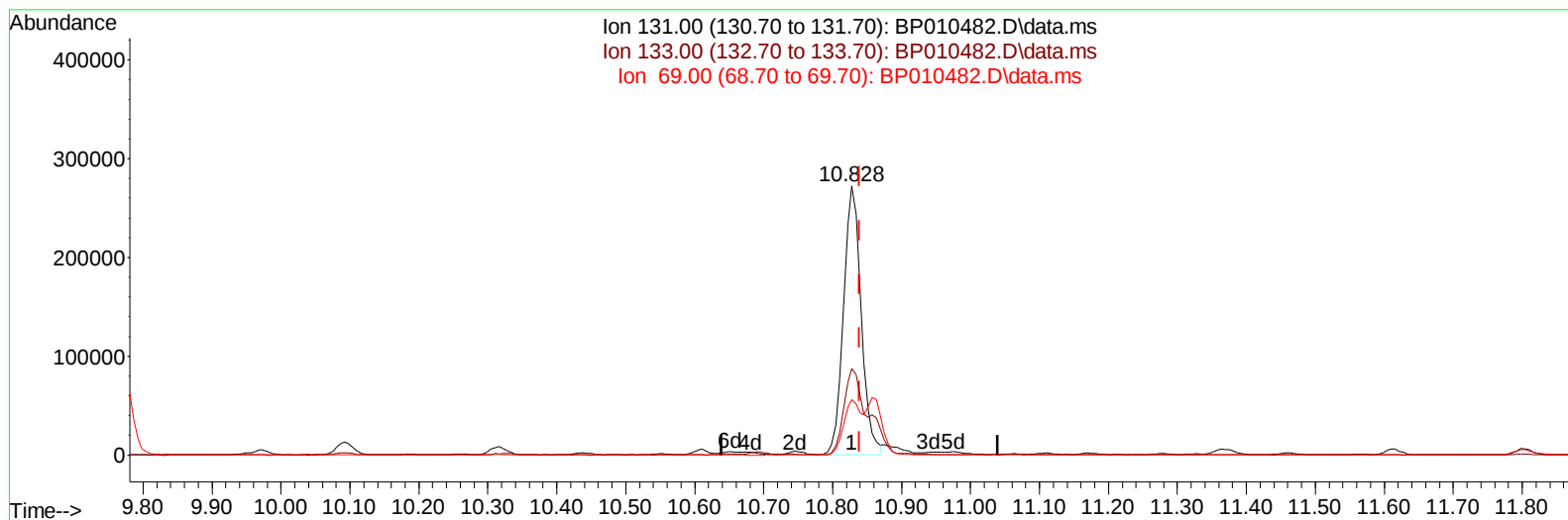
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052322\
 Data File : BP010482.D
 Acq On : 23 May 2022 18:18
 Operator : CG/JU
 Sample : N2918-18
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTD4

Manual IntegrationsAPPROVED

Quant Time: May 23 23:20:18 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 13 15:04:31 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/24/2022
 Supervised By :Yogesh Patel 05/26/2022



TIC: BP010482.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.828min (-0.012) 32.25 ng/ul

response 483088

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	32.18
69.00	26.10	20.61#
0.00	0.00	0.00

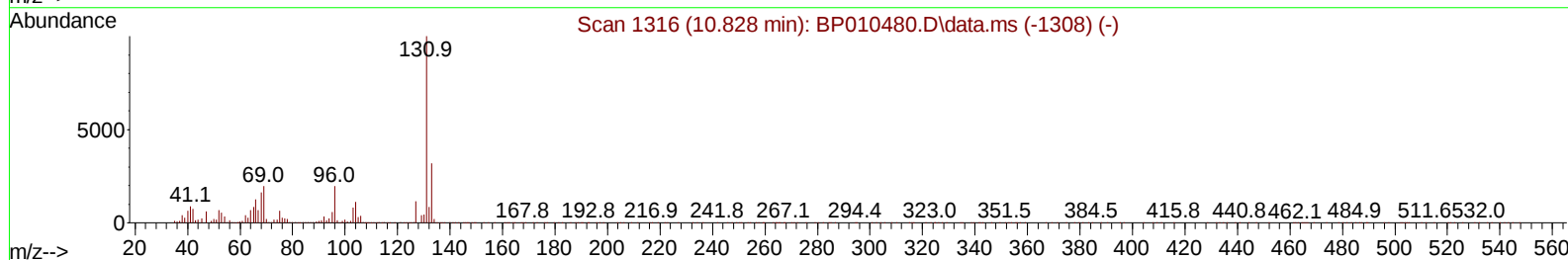
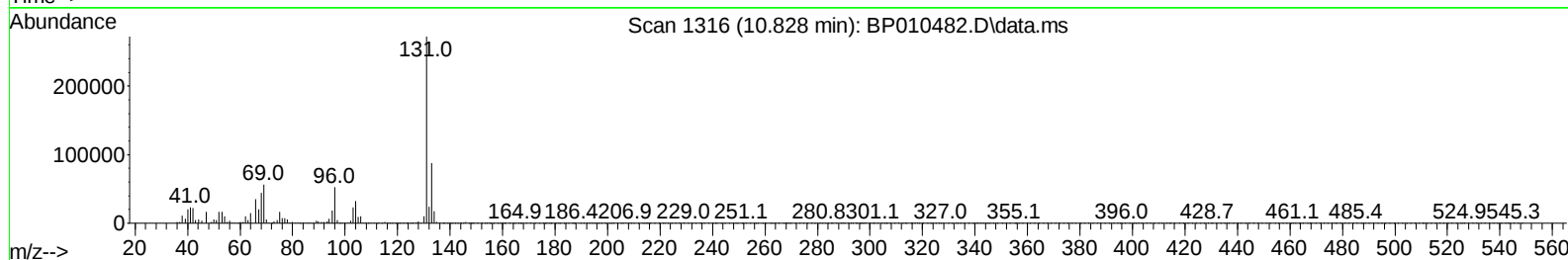
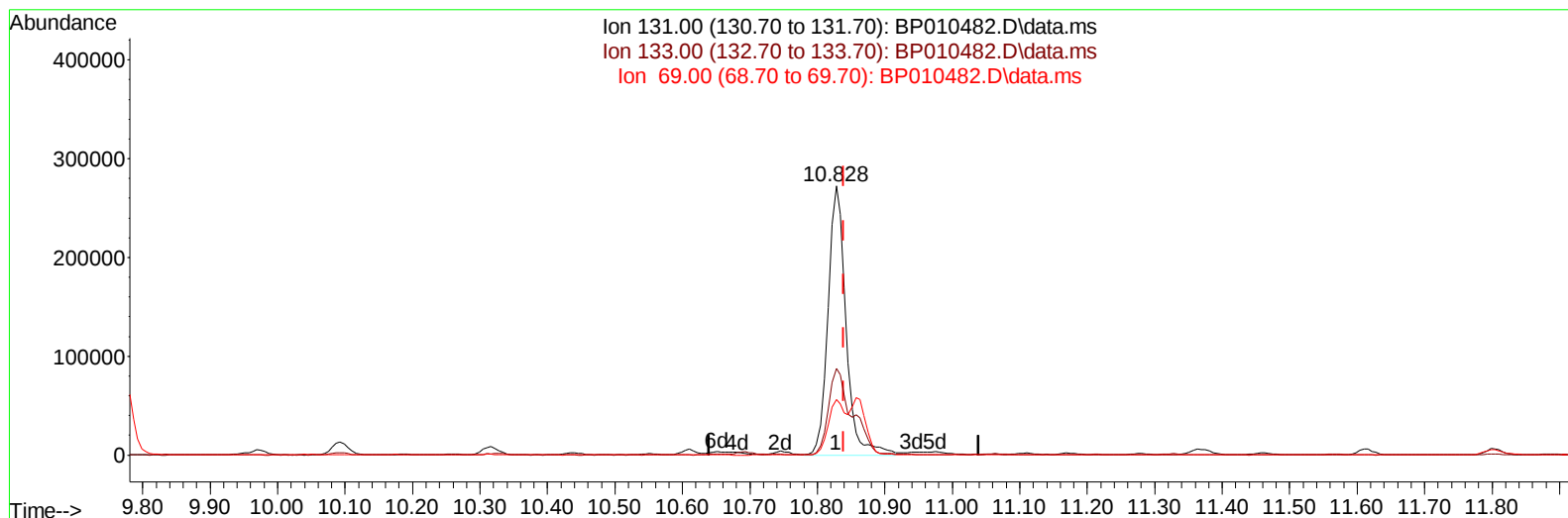
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TIC: BP010482.D\data.ms

(31) 4-Chloroaniline-d4 (S)

10.828min (-0.012) 33.67 ng/ul m

response 504465

Ion	Exp%	Act%
131.00	100.00	100.00
133.00	29.20	32.18
69.00	26.10	20.61#
0.00	0.00	0.00

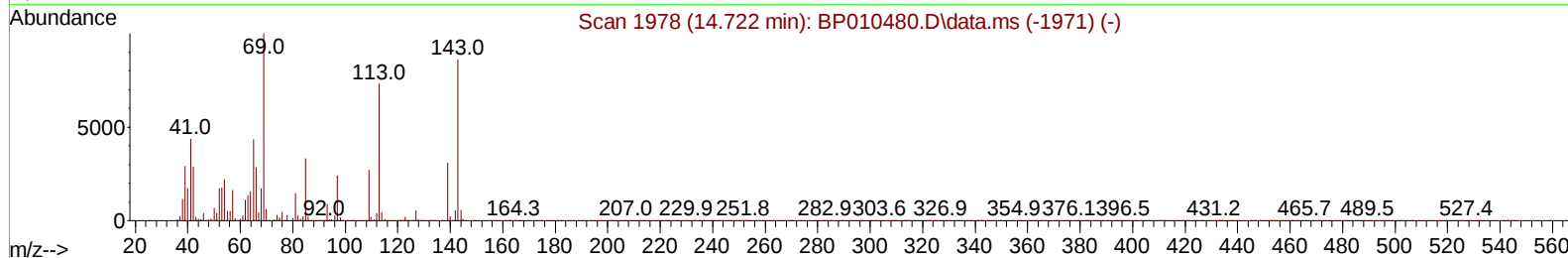
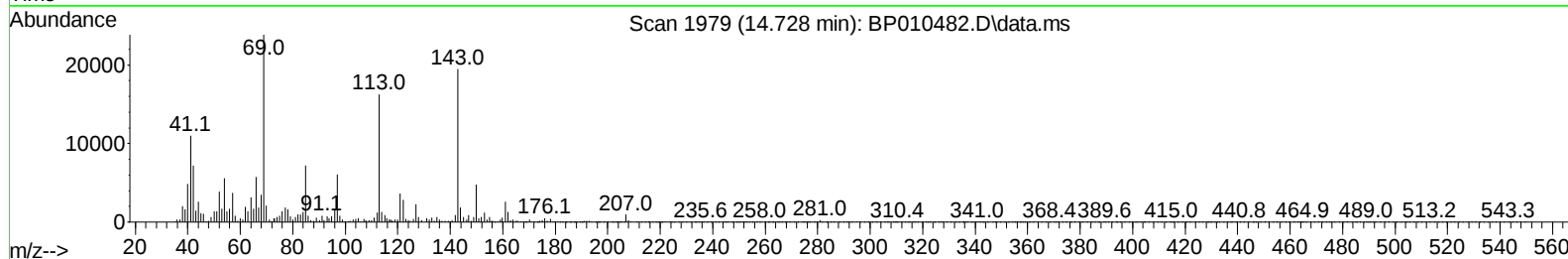
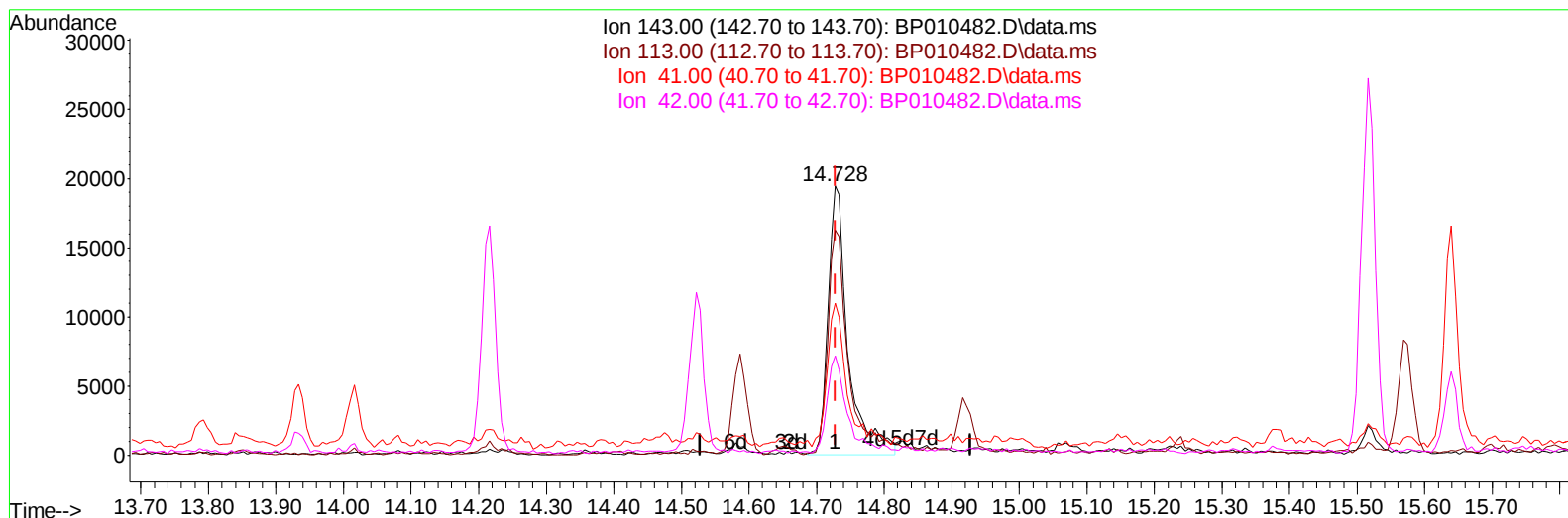
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Instrument :
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TIC: BP010482.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (+ 0.000) 7.05 ng/ul m

response 40234

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	83.53#
41.00	23.20	56.40#
42.00	18.00	36.91#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.893	152	145749	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	649672	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	435109	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	944242	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	851475	20.000	ng/ul	# 0.00
88) Perylene-d12	23.774	264	635208	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	18236	5.333	ng/uL	0.00
4) Pyridine-d5	3.764	84	100294	10.330	ng/ul	0.00
7) Phenol-d5	7.057	99	102593	8.387	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	337822	41.103	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	299150	31.958	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	212206	20.412	ng/ul	-0.01
21) Nitrobenzene-d5	9.051	128	217585	43.274	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	225721	42.361	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.316	165	371047	36.827	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	504465m	33.674	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	1428709	44.567	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1622642	43.882	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	40234m	7.051	ng/ul	0.00
60) Fluorene-d10	15.516	176	1235855	44.313	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	209417	37.922	ng/ul	0.00
73) Anthracene-d10	17.375	188	1991693	46.582	ng/ul	0.00
81) Pyrene-d10	19.604	212	2359870	52.187	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.621	264	1562268	48.980	ng/ul	-0.01
Target Compounds						
					Qvalue	
30) Naphthalene	10.746	128	3047078	89.796	ng/ul	97
37) 1-Methylnaphthalene	12.569	142	72027	2.970	ng/ul#	97
52) Acenaphthene	14.586	153	1267773	48.582	ng/ul	97
56) Dibenzofuran	14.922	168	103895	2.717	ng/ul	100
61) Fluorene	15.569	166	528596	16.902	ng/ul	92
72) Phenanthrene	17.316	178	74935	1.487	ng/ul	98
77) Carbazole	17.680	167	1735870	39.071	ng/ul#	96

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

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