

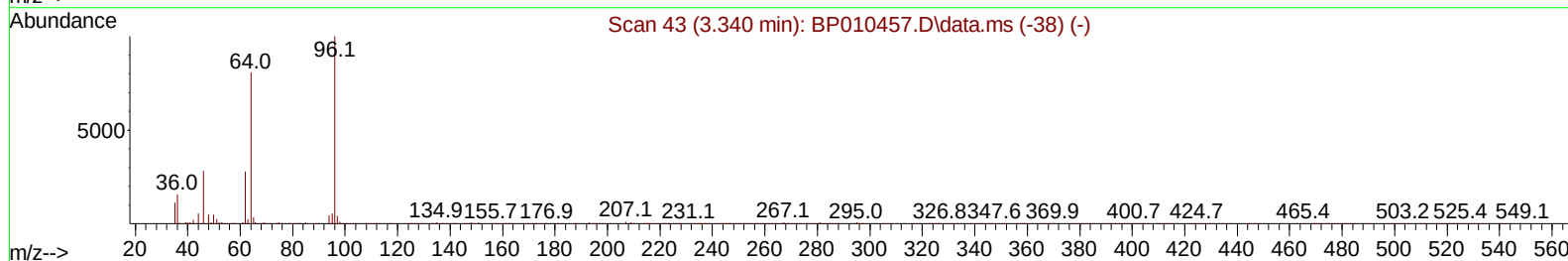
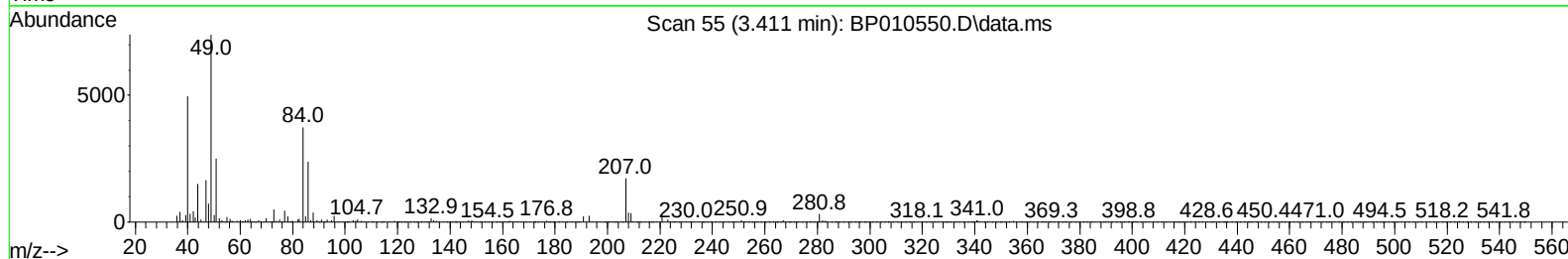
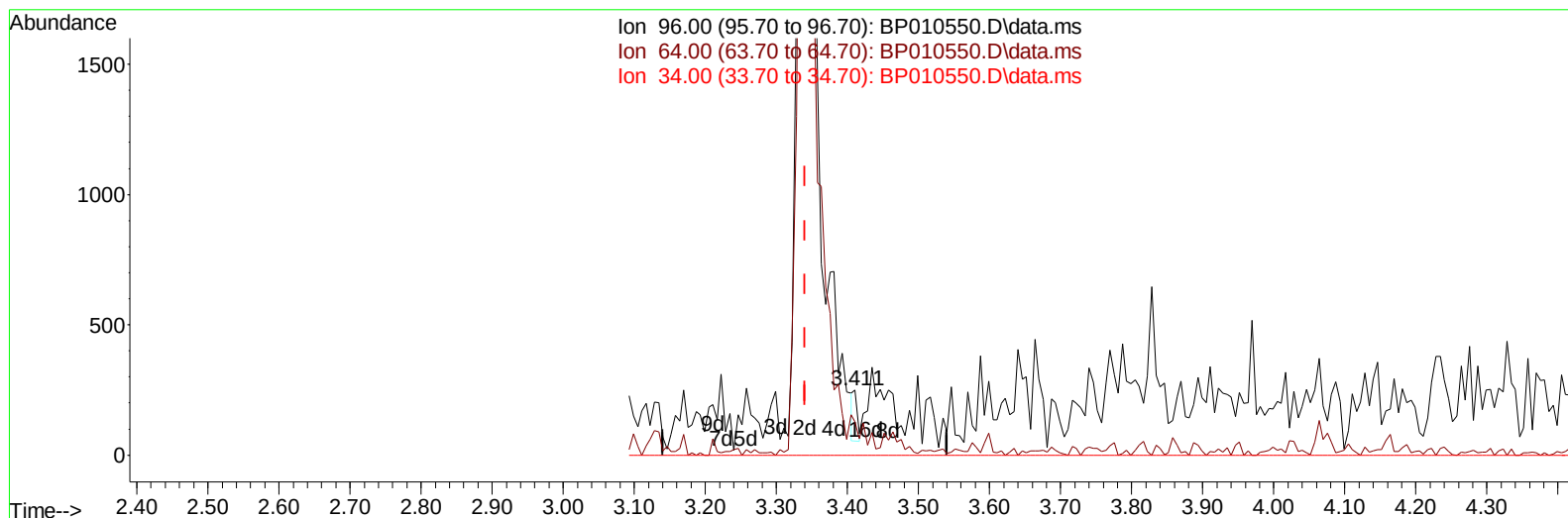
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
 Data File : BP010550.D
 Acq On : 25 May 2022 23:13
 Operator : CG/JU
 Sample : N2883-03
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4K5

Manual IntegrationsAPPROVED

Quant Time: May 26 00:54:34 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 24 05:50:40 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/26/2022
 Supervised By :mohammad ahmed 05/27/2022



TIC: BP010550.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.411min (+ 0.071) 0.03 ng/uL

response 79

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	59.04
34.00	0.00	0.00
0.00	0.00	0.00

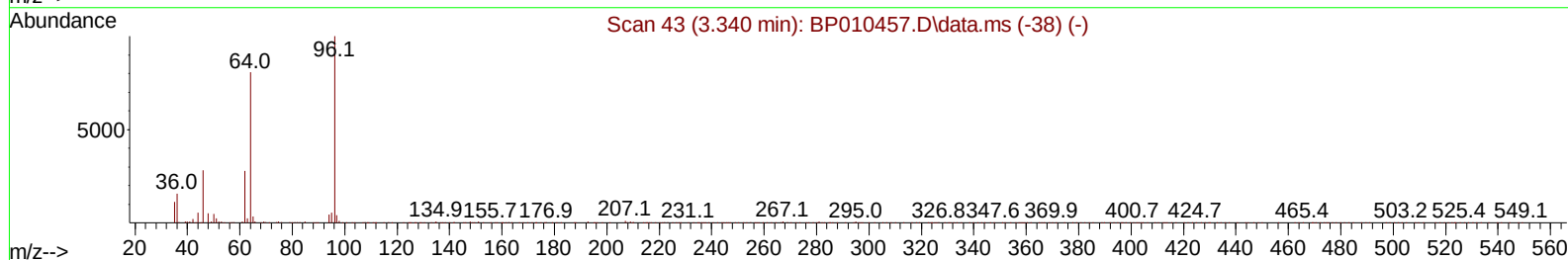
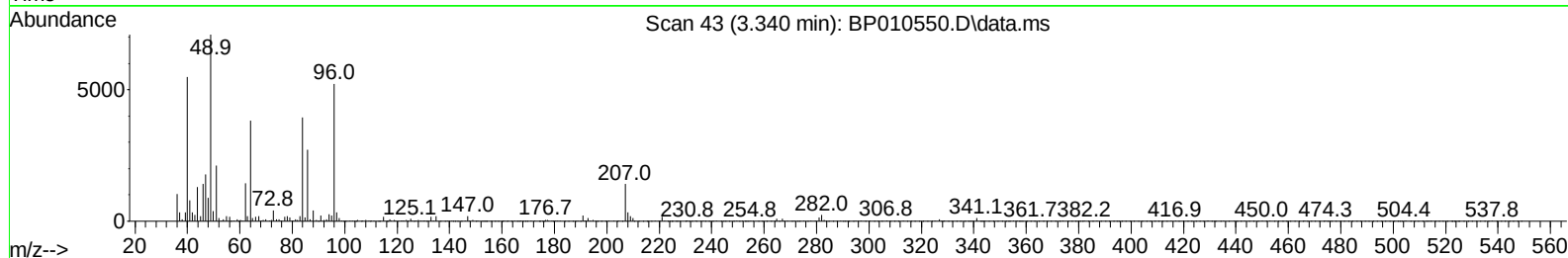
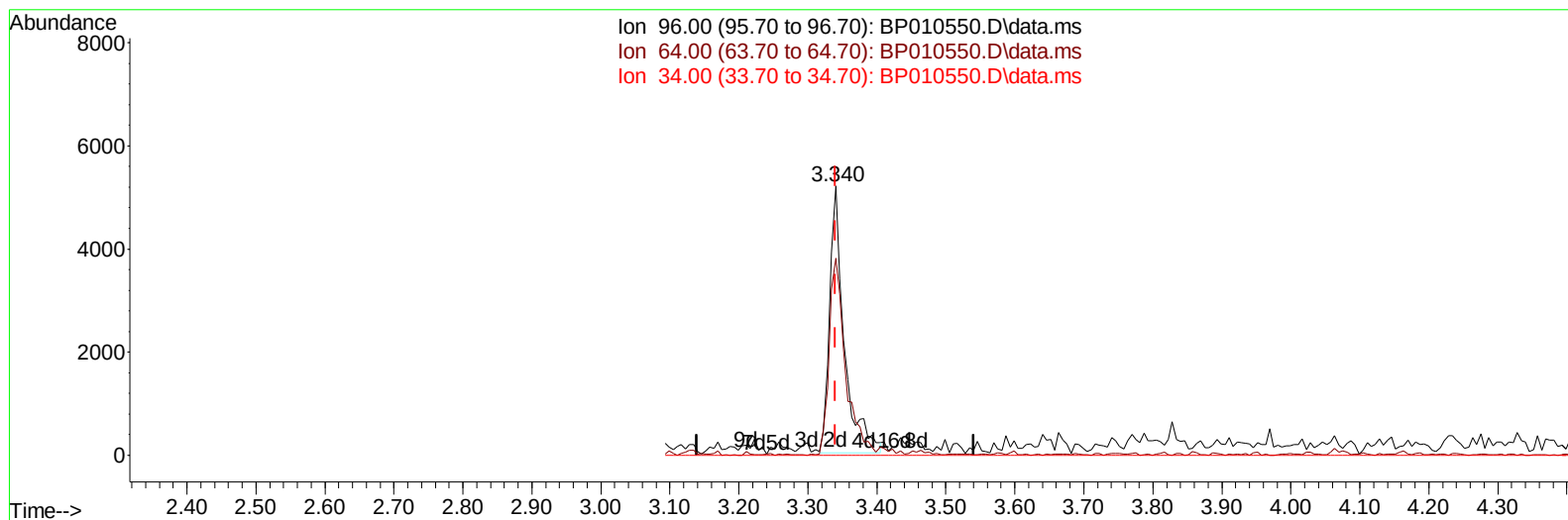
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TIC: BP010550.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.340min (0.000) 2.98 ng/uL m

response 7674

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	73.14#
34.00	0.00	0.00
0.00	0.00	0.00

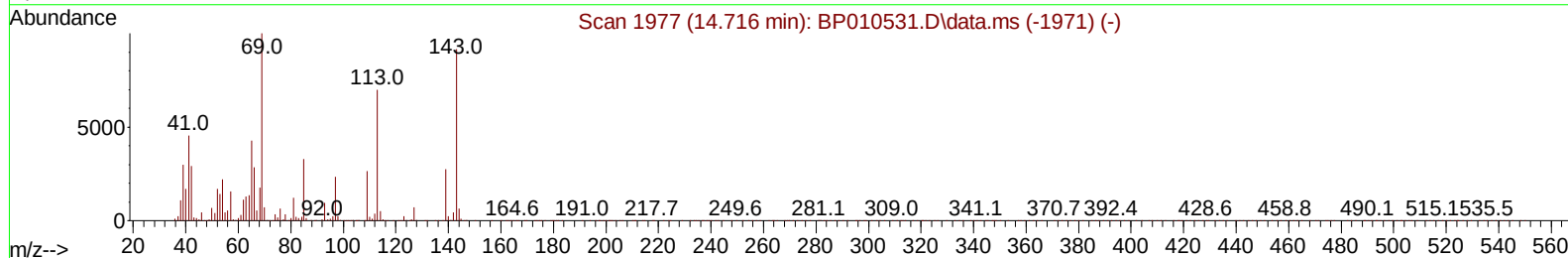
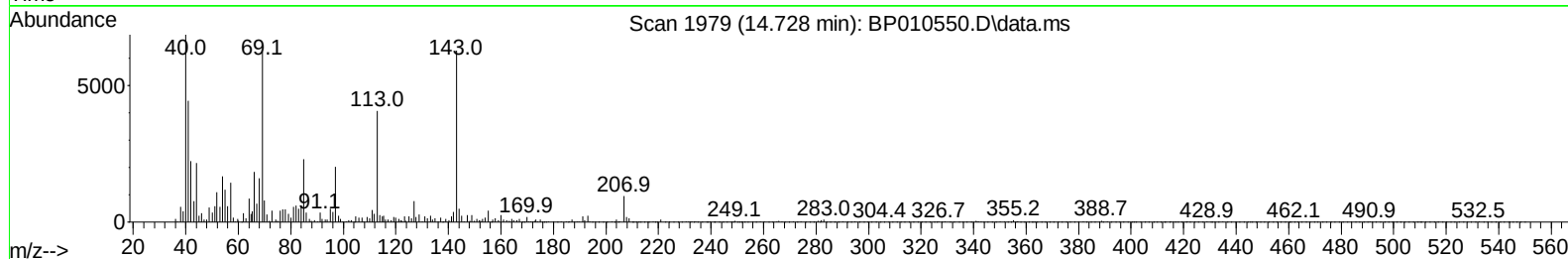
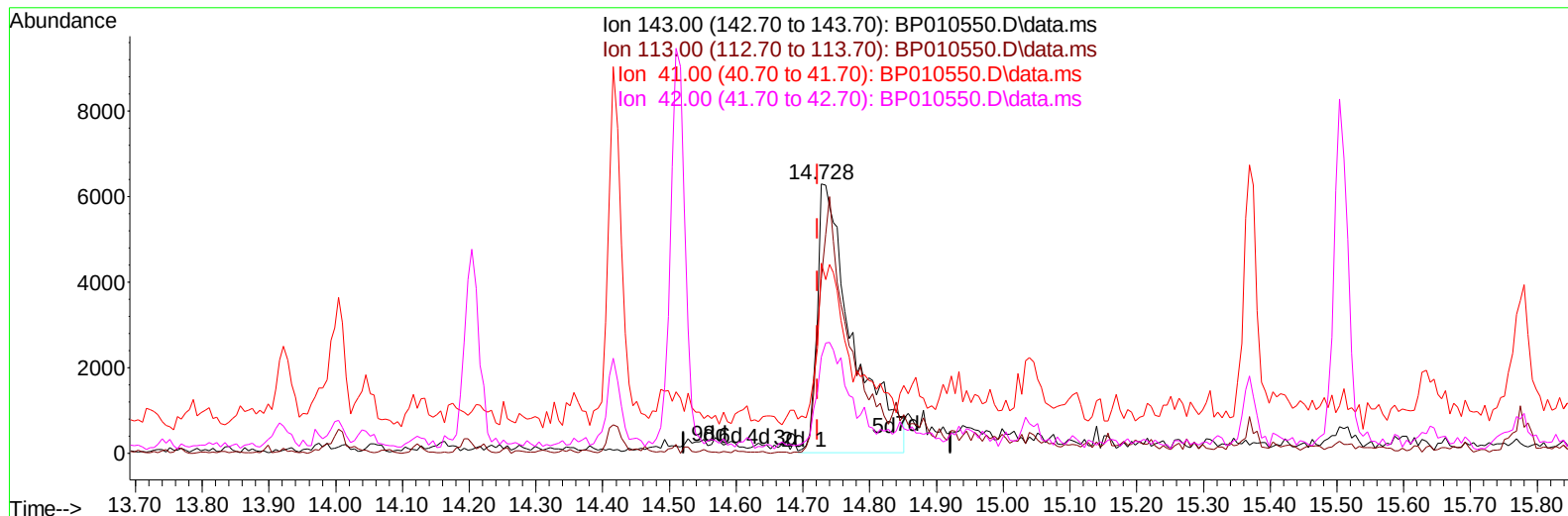
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ALS Vial : 20 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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Manual IntegrationsAPPROVED

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QLast Update : Tue May 24 05:50:40 2022
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TIC: BP010550.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.728min (+ 0.006) 5.22 ng/ul m

response 23049

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	64.62#
41.00	23.20	70.69#
42.00	18.00	35.66#

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 Misc :
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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	7.881	152	109586	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	470852	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	336907	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.263	188	749510	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.345	240	741741	20.000	ng/ul	0.00
88) Perylene-d12	23.763	264	682772	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	7674m	2.985	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.052	99	120331	13.083	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.210	67	97826	15.830	ng/ul	0.00
11) 2-Chlorophenol-d4	7.416	132	104996	14.918	ng/ul	0.00
15) 4-Methylphenol-d8	8.593	113	103344	13.221	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	53617	14.713	ng/ul	0.00
24) 2-Nitrophenol-d4	9.769	143	51633	13.370	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	89483	12.254	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	93230	8.587	ng/ul	0.00
46) Dimethylphthalate-d6	13.922	166	371214	14.955	ng/ul	0.00
49) Acenaphthylene-d8	14.204	160	430401	15.032	ng/ul	0.00
54) 4-Nitrophenol-d4	14.728	143	23049m	5.217	ng/ul	0.00
60) Fluorene-d10	15.504	176	333692	15.452	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	17289	3.944	ng/ul	0.00
73) Anthracene-d10	17.363	188	538012	15.852	ng/ul	0.00
81) Pyrene-d10	19.598	212	757493	19.230	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.604	264	609773	17.786	ng/ul	-0.01
Target Compounds						
					Qvalue	
86) Bis(2-ethylhexyl)phtha...	21.257	149	41121	1.467	ng/ul#	82

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

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