

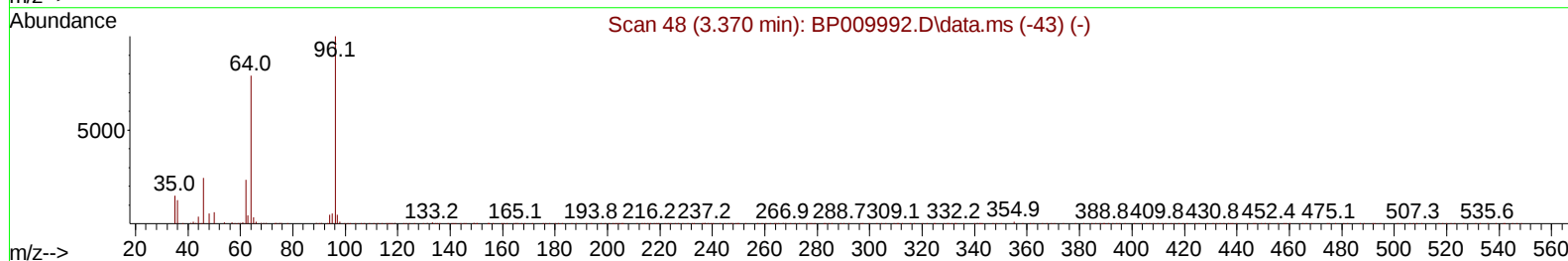
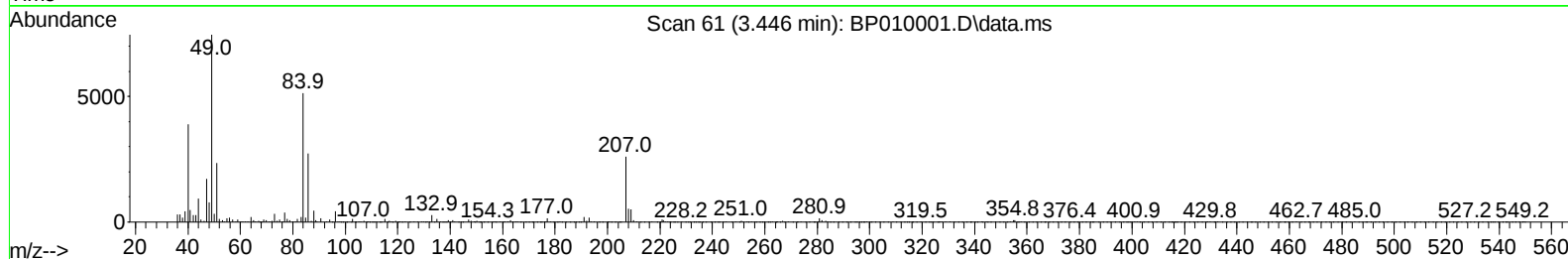
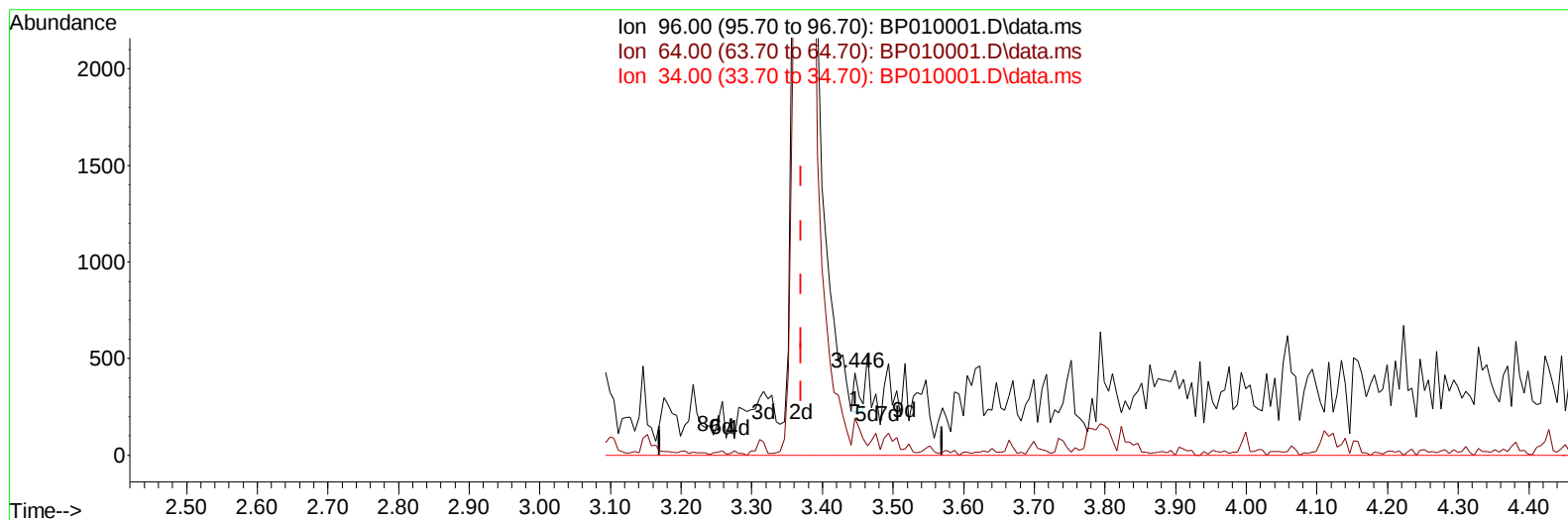
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041922\
 Data File : BP010001.D
 Acq On : 21 Apr 2022 05:37
 Operator : CG/JU
 Sample : N2472-14
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J59

Manual IntegrationsAPPROVED

Quant Time: Apr 21 06:54:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 20 01:32:14 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/21/2022
 Supervised By :Yogesh Patel 04/25/2022



TIC: BP010001.D\data.ms

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

3.446min (+ 0.076) 0.04 ng/uL

response 109

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.07
34.00	0.00	0.00
0.00	0.00	0.00

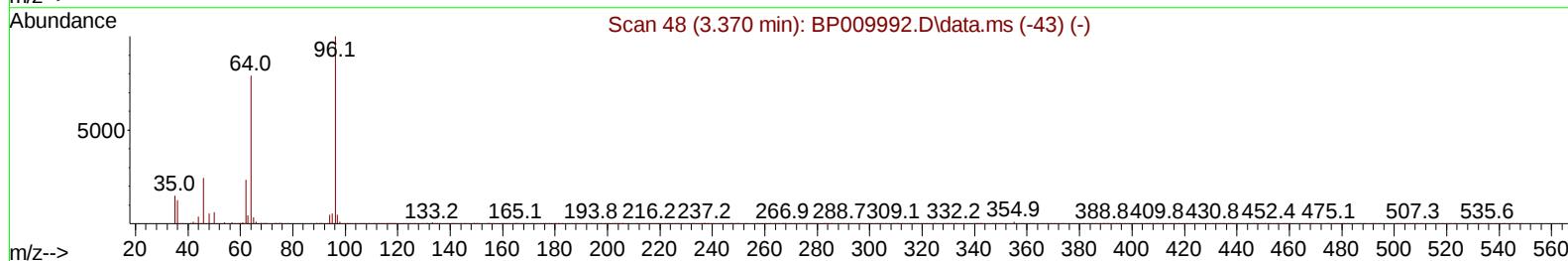
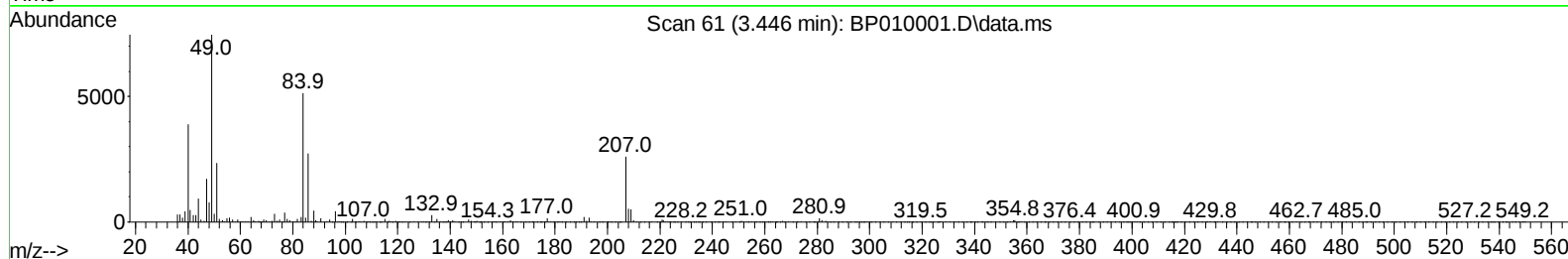
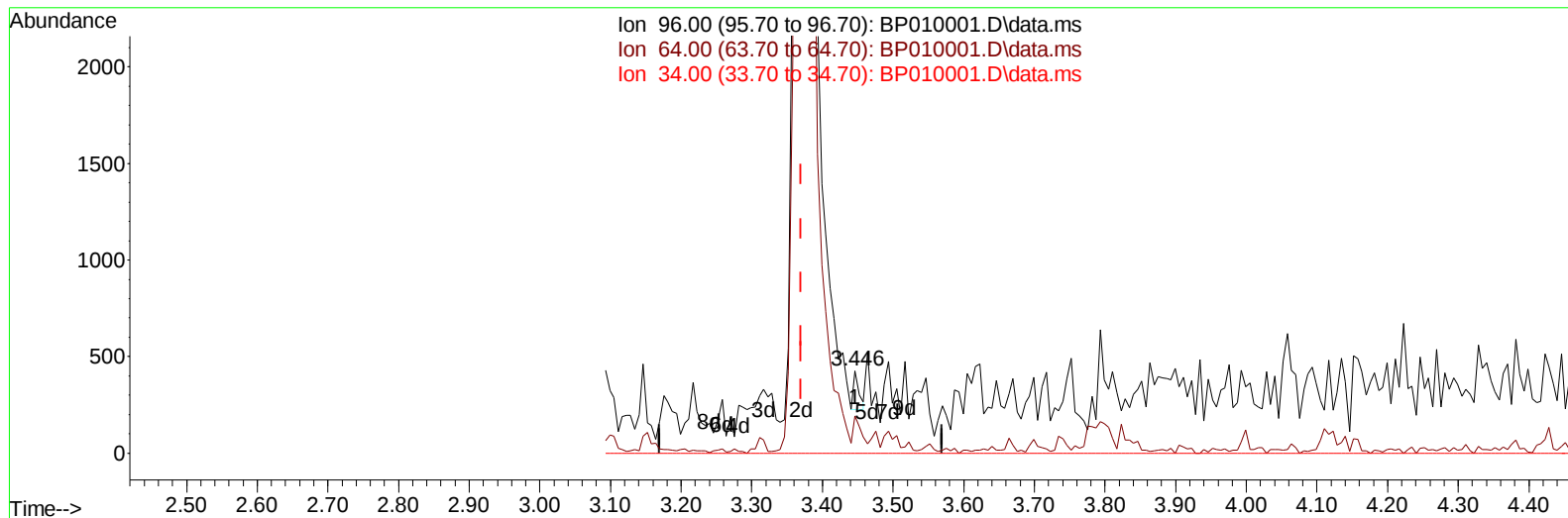
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(3) 1,4-Dioxane-d8 (S)

3.446min (+ 0.076) 0.04 ng/uL

response 109

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	45.07
34.00	0.00	0.00
0.00	0.00	0.00

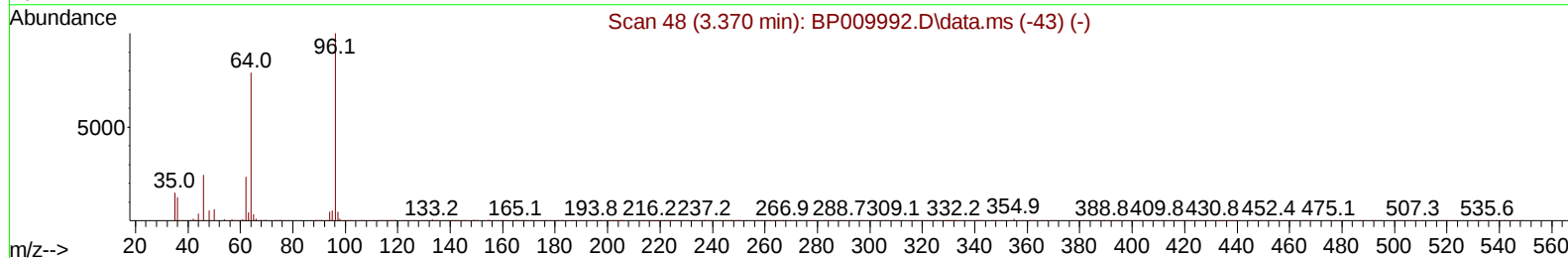
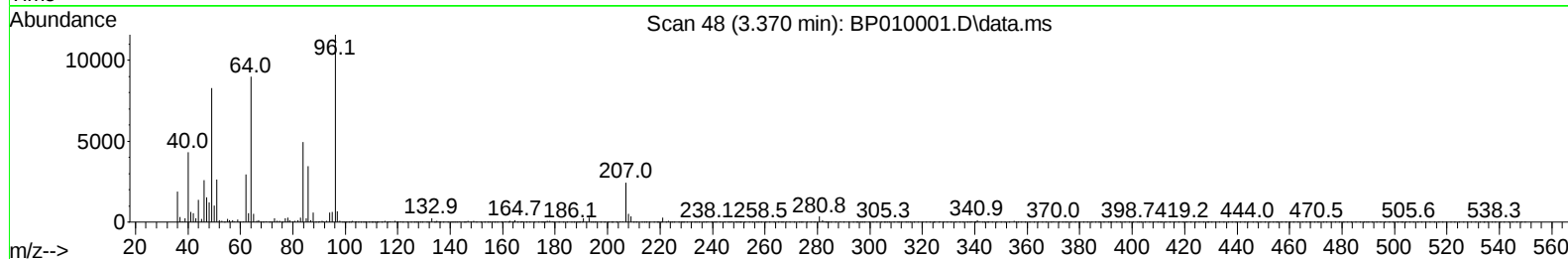
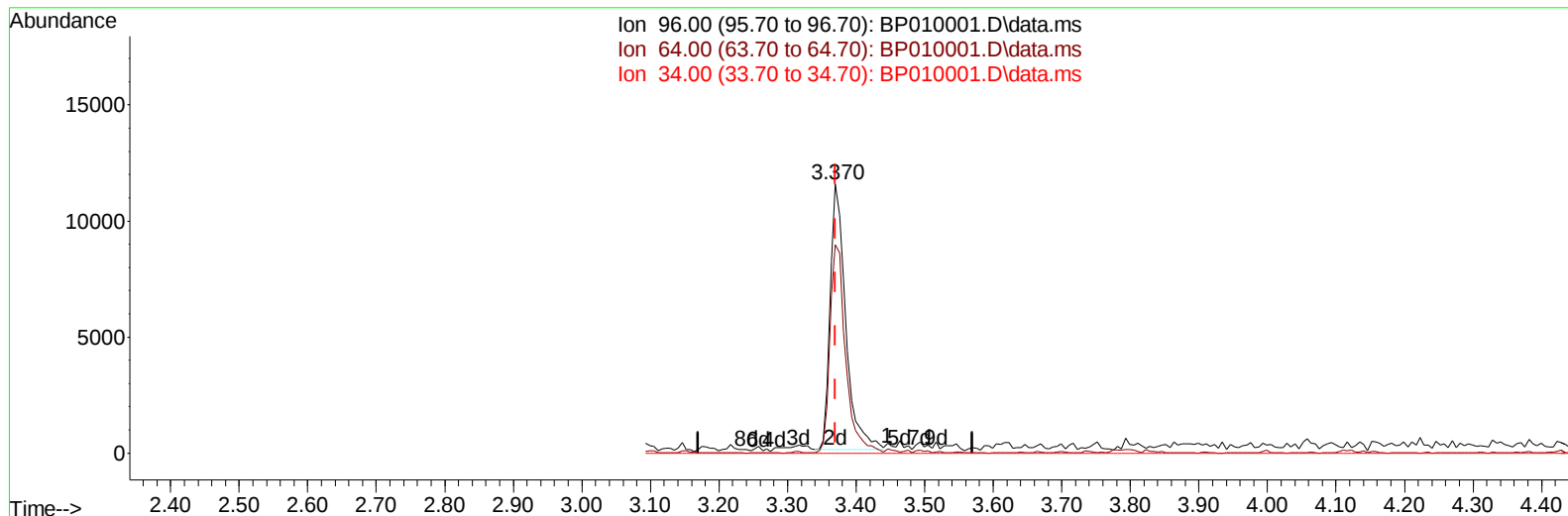
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TIC: BP010001.D\data.ms

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

3.370min (+ 0.000) 6.03 ng/uL m

response 17992

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	77.56#
34.00	0.00	0.00
0.00	0.00	0.00

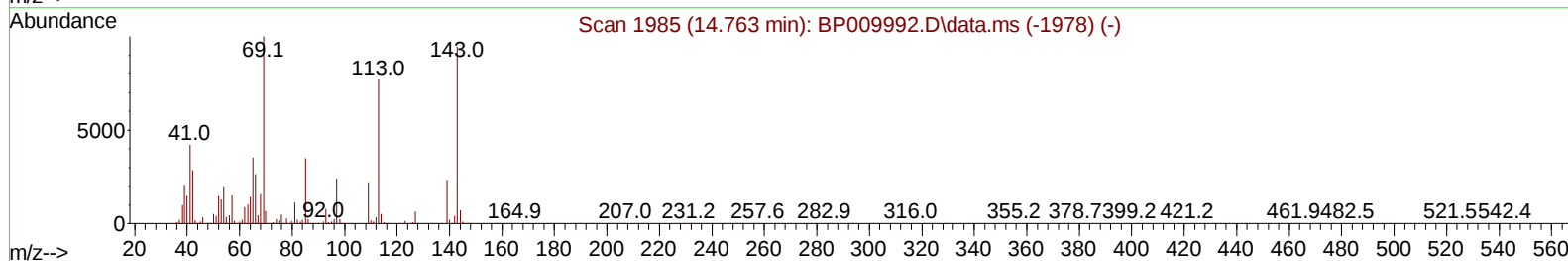
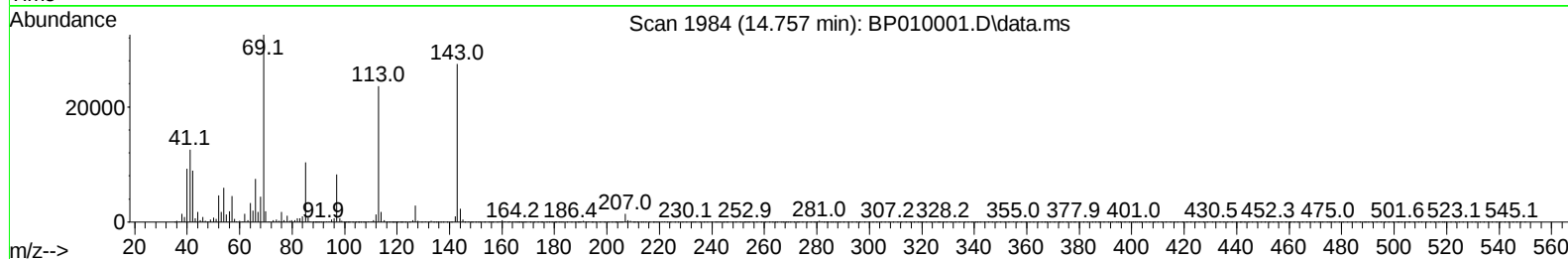
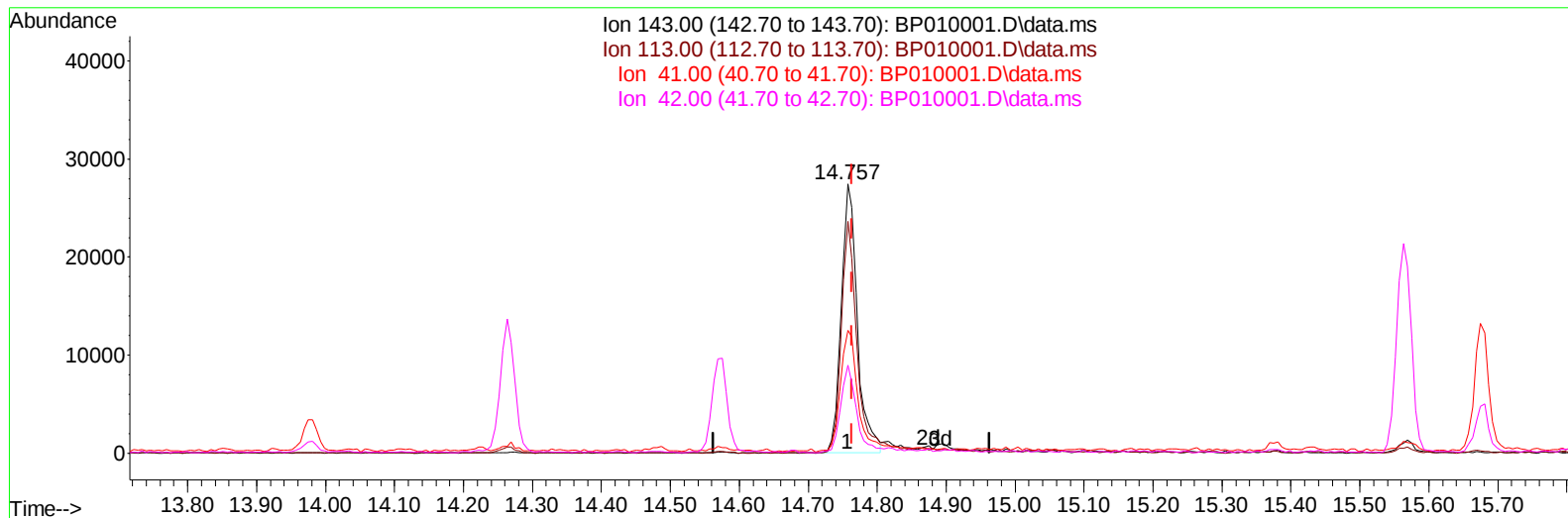
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TIC: BP010001.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 7.17 ng/u1

response 44287

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.13#
41.00	23.20	45.71#
42.00	18.00	32.73#

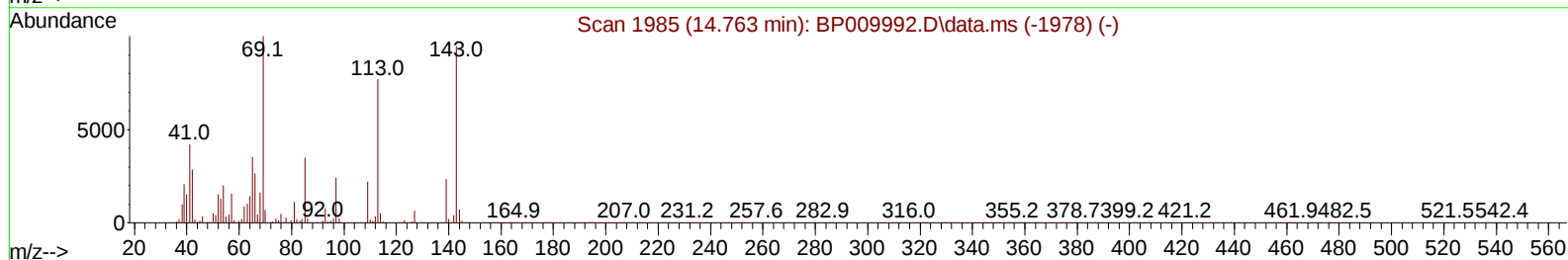
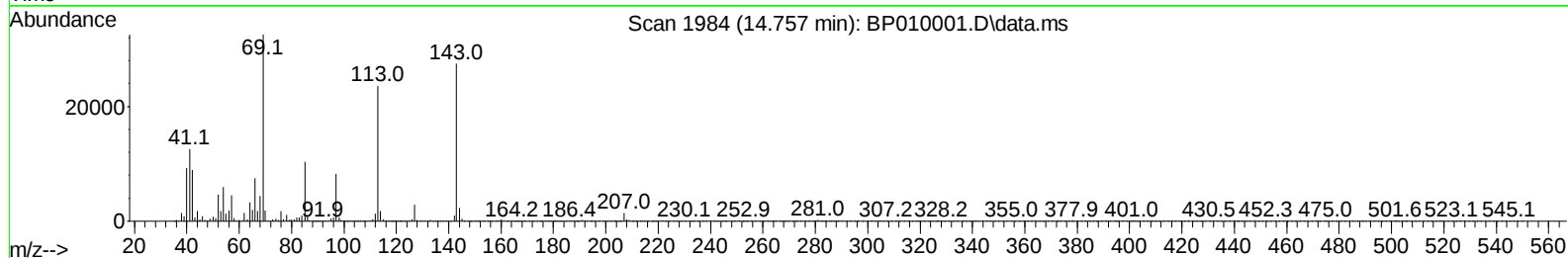
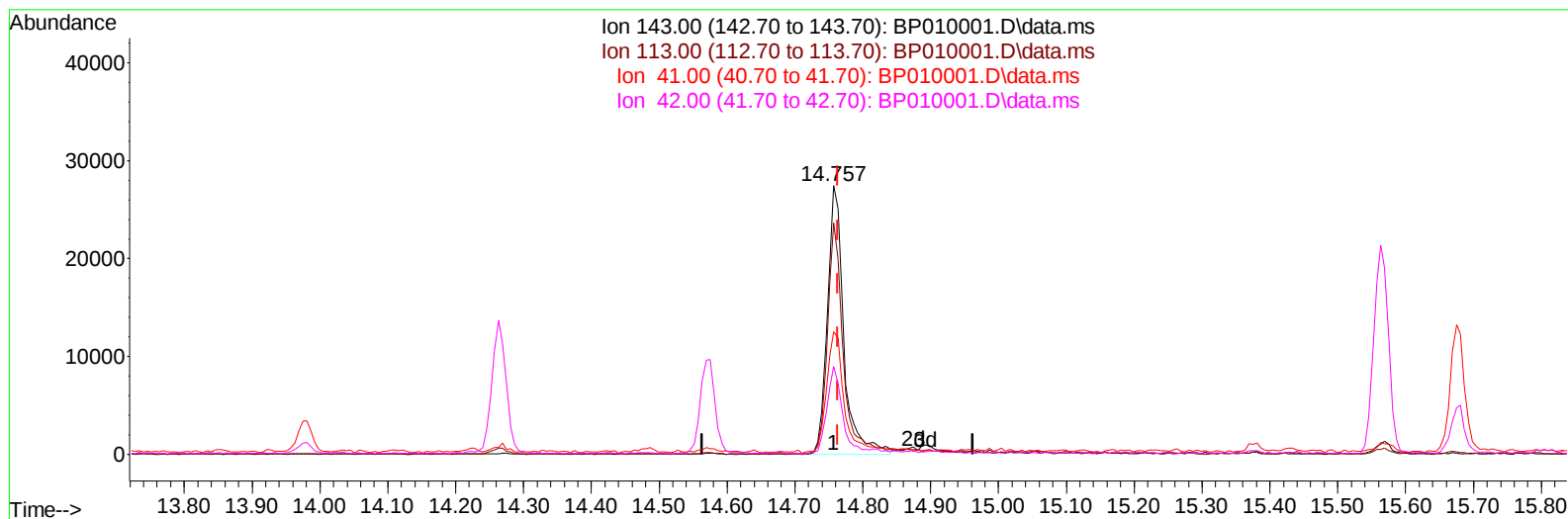
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Operator : CG/JU
Sample : N2472-14
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0J59

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/21/2022
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TIC: BP010001.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.757min (-0.006) 7.47 ng/ul m

response 46178

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	86.13#
41.00	23.20	45.71#
42.00	18.00	32.73#

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 Sample : N2472-14
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	133212	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	601942	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	433859	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	956347	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	866697	20.000	ng/ul	# 0.00
88) Perylene-d12	23.851	264	728901	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	17992m	6.028	ng/uL	0.00
4) Pyridine-d5	3.787	84	94978	11.541	ng/ul	0.00
7) Phenol-d5	7.099	99	99083	8.472	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	267013	38.350	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	268144	30.395	ng/ul	0.00
15) 4-Methylphenol-d8	8.646	113	195286	20.003	ng/ul	0.00
21) Nitrobenzene-d5	9.099	128	173666	40.008	ng/ul	0.00
24) 2-Nitrophenol-d4	9.828	143	187636	39.427	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	334983	33.638	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131	419396	29.569	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166	1319044	38.704	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1464774	36.279	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	46178m	7.471	ng/ul	0.00
60) Fluorene-d10	15.563	176	1170474	40.826	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	206591	36.465	ng/ul	-0.01
73) Anthracene-d10	17.422	188	1919845	41.928	ng/ul	0.00
81) Pyrene-d10	19.651	212	2324605	49.110	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1668945	43.968	ng/ul	-0.01

Target Compounds Qvalue

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

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