

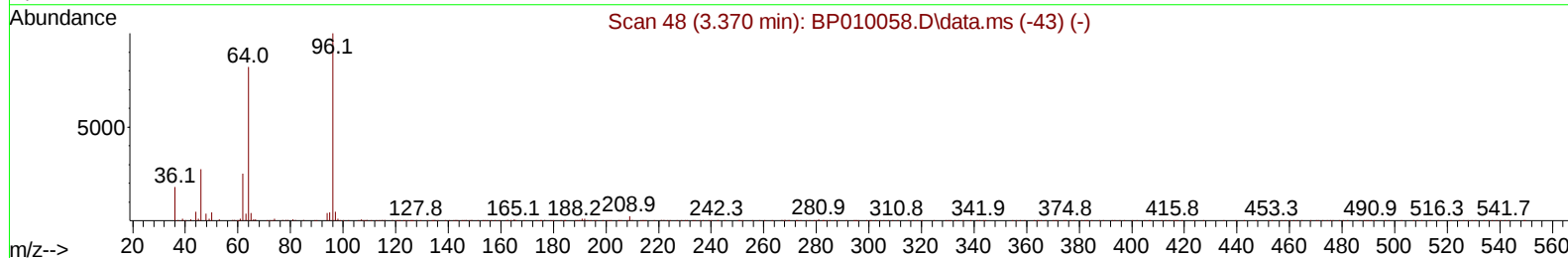
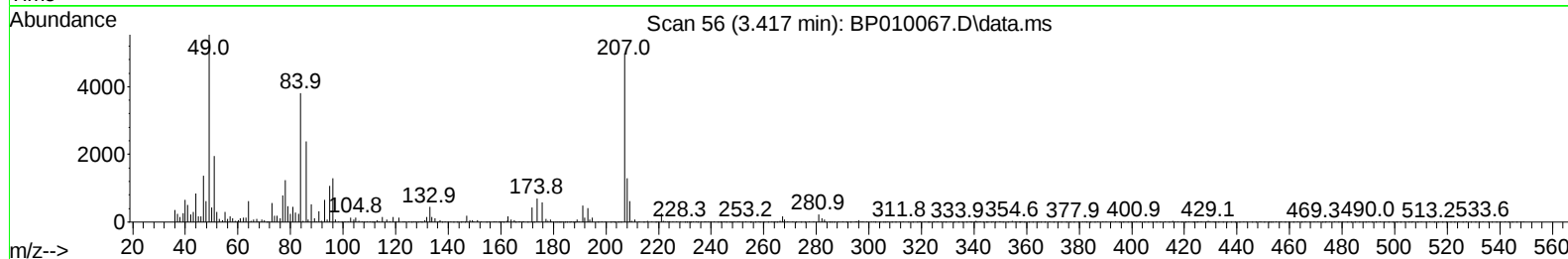
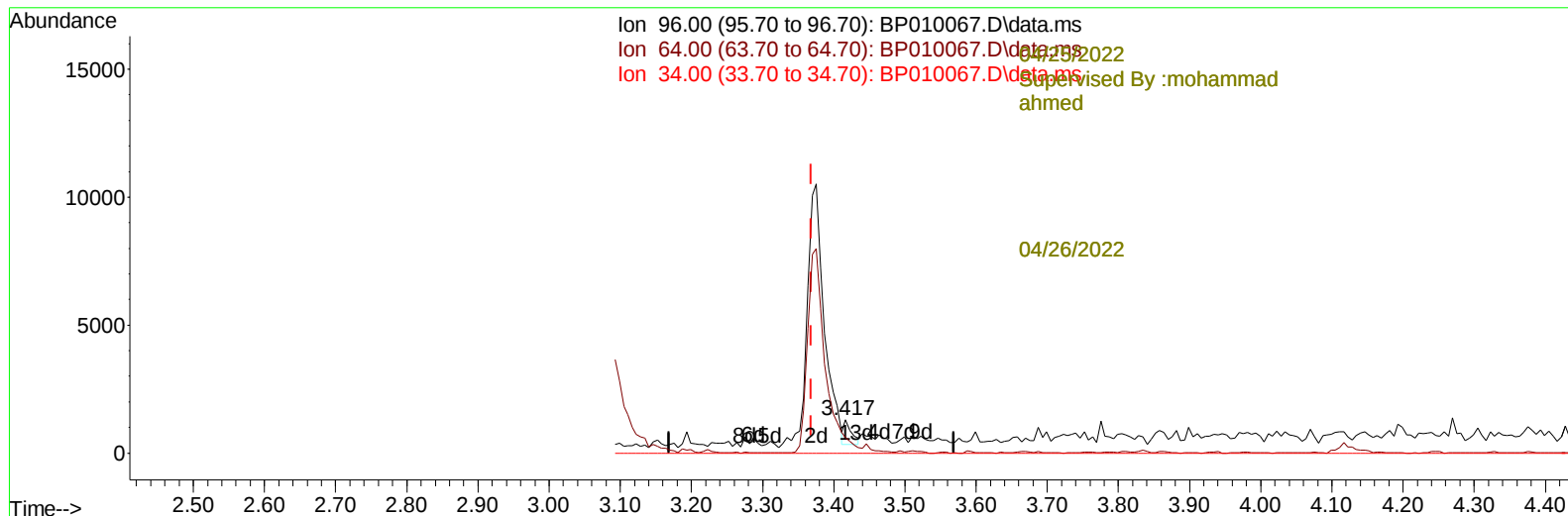
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
 Data File : BP010067.D
 Acq On : 22 Apr 2022 22:14
 Operator : CG/JU
 Sample : N2490-07
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J44

Manual IntegrationsAPPROVED

Quant Time: Apr 22 23:32:35 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 21 14:49:38 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/25/2022
 Supervised By :mohammad ahmed 04/26/2022



TIC: BP010067.D\data.ms

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

3.417min (+ 0.048) 0.22 ng/uL

response 726

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	47.19
34.00	0.00	0.00
0.00	0.00	0.00

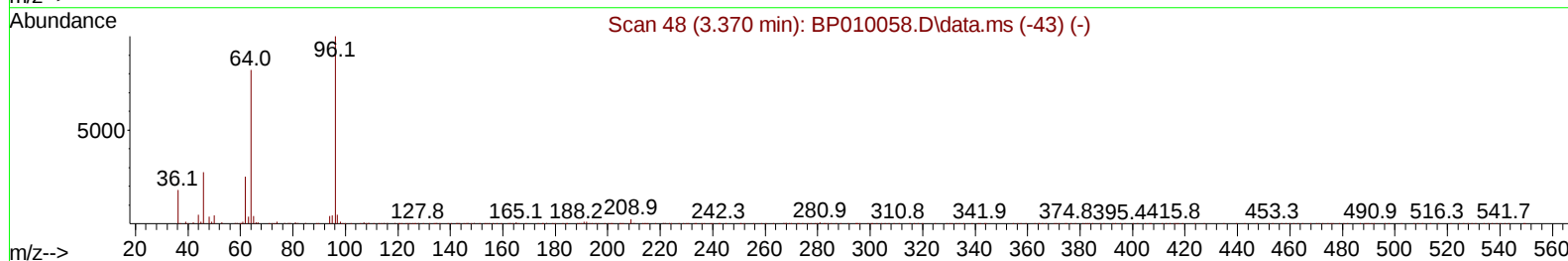
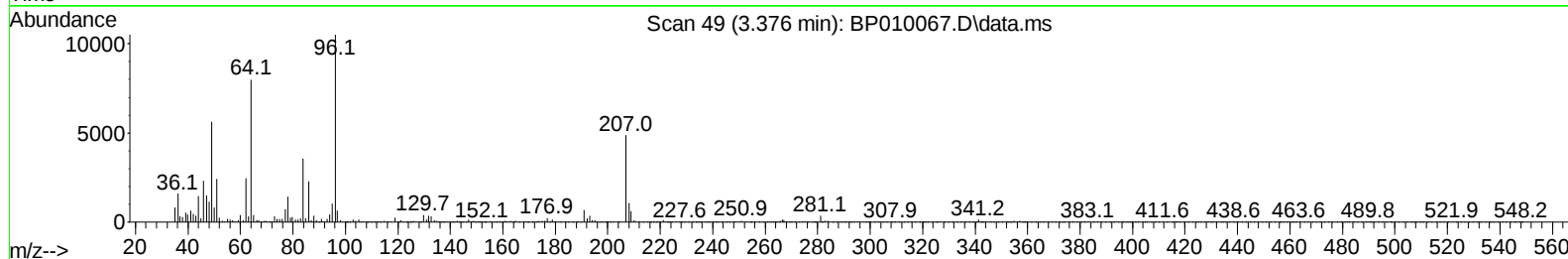
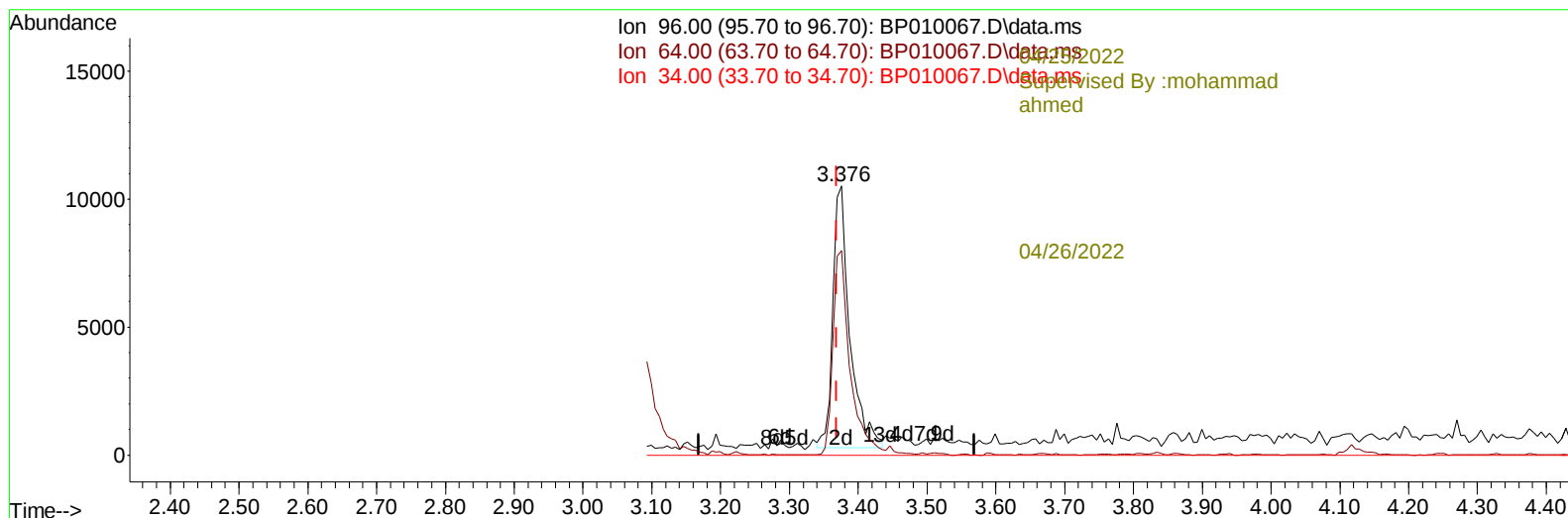
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(3) 1,4-Dioxane-d8 (S)

3.376min (+ 0.006) 5.34 ng/uL m

response 17546

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	75.93#
34.00	0.00	0.00
0.00	0.00	0.00

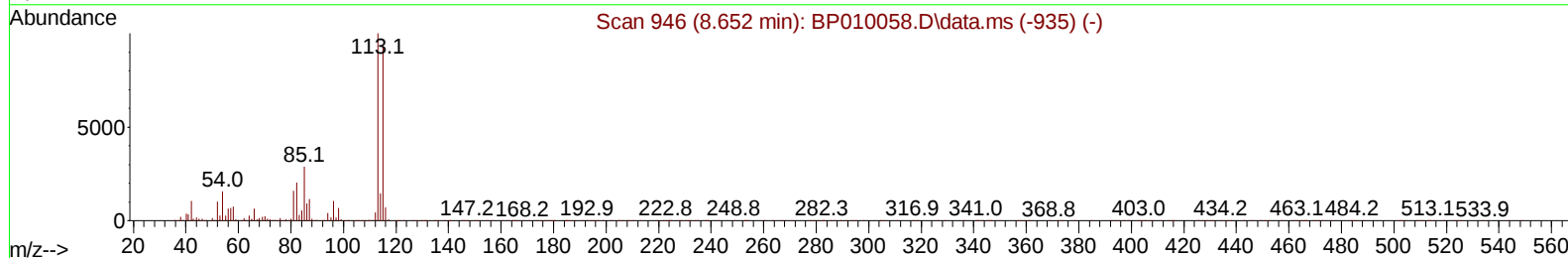
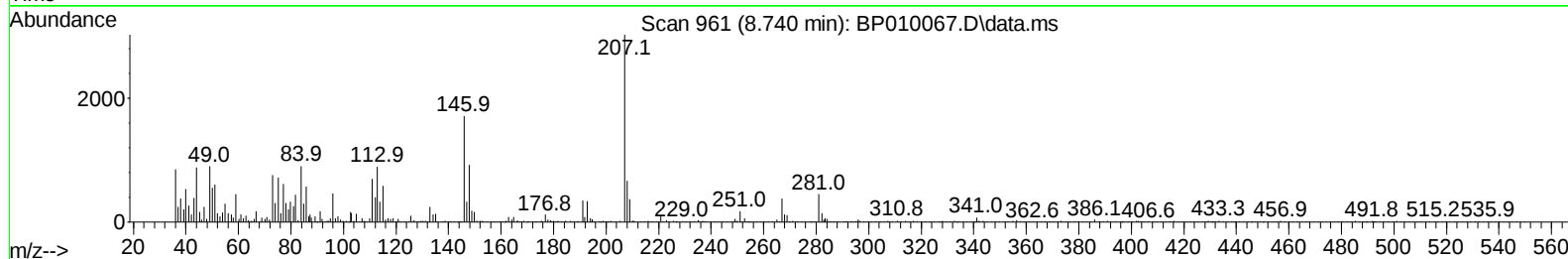
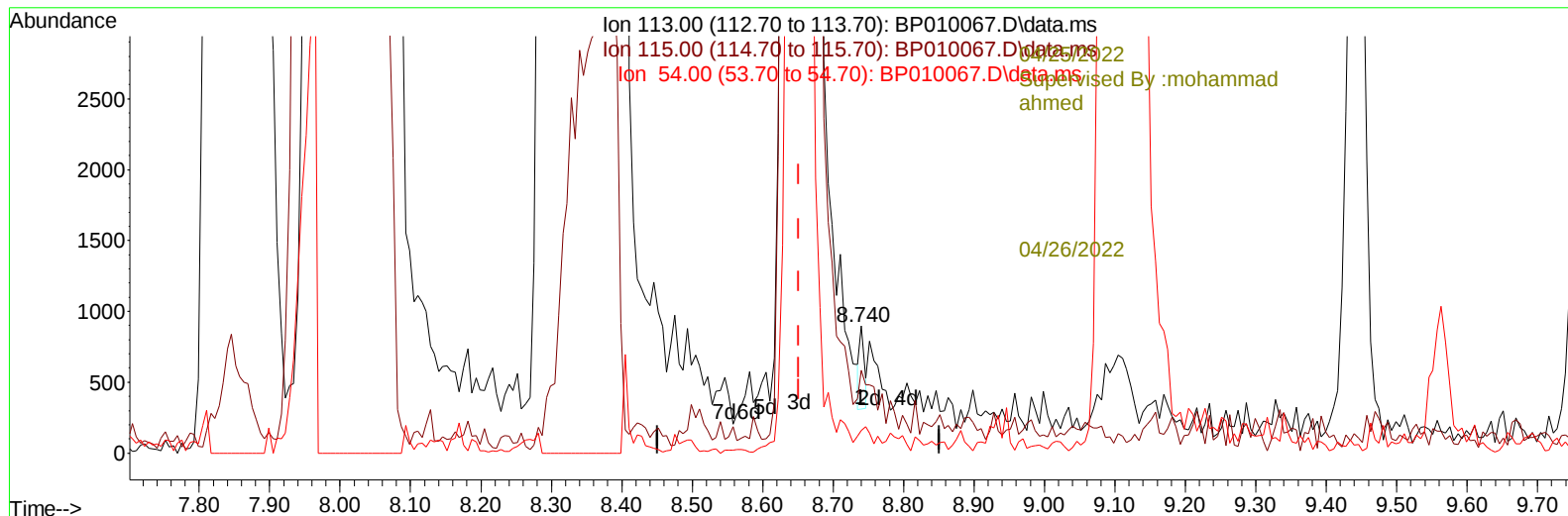
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Instrument :
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TIC: BP010067.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.740min (+ 0.088) 0.03 ng/u1

response 283

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	106.70	65.11#
54.00	4.50	17.73#
0.00	0.00	0.00

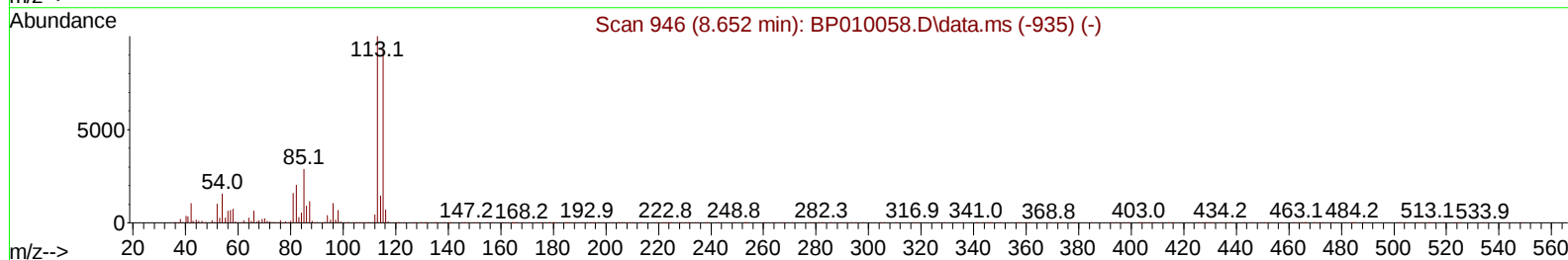
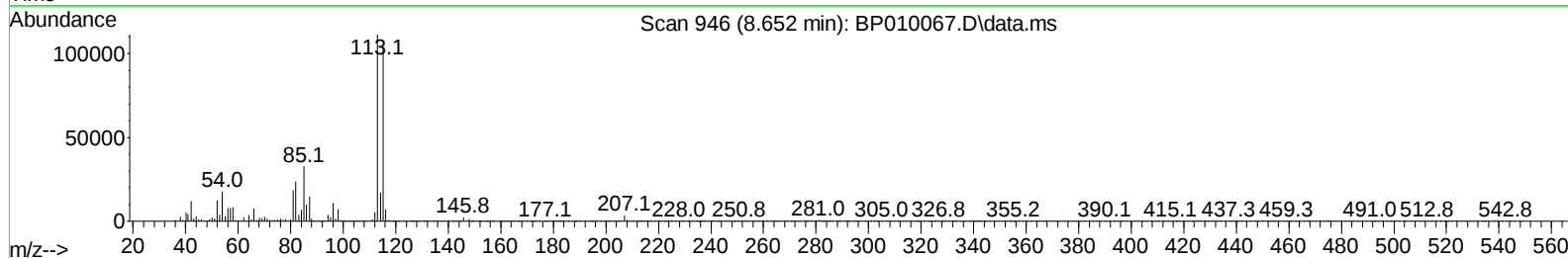
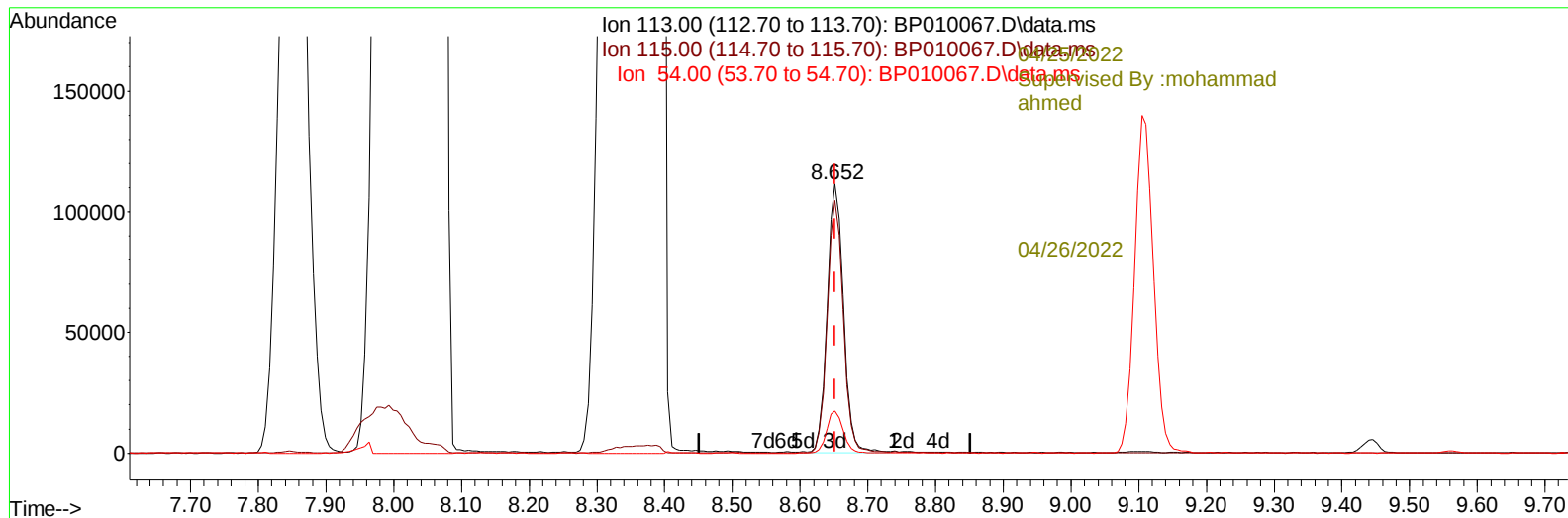
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TIC: BP010067.D\data.ms

(15) 4-Methylphenol-d8 (S)

8.652min (+ 0.000) 17.18 ng/ul m

response 185634

Ion	Exp%	Act%
113.00	100.00	100.00
115.00	106.70	94.06
54.00	4.50	15.80#
0.00	0.00	0.00

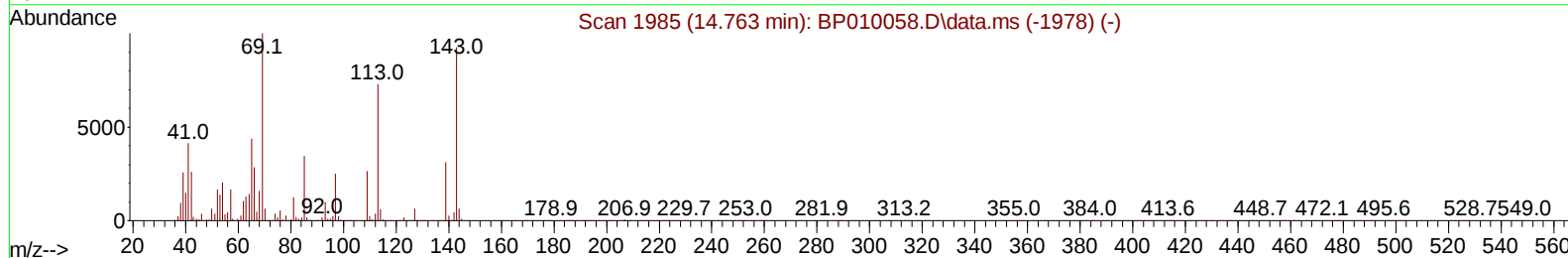
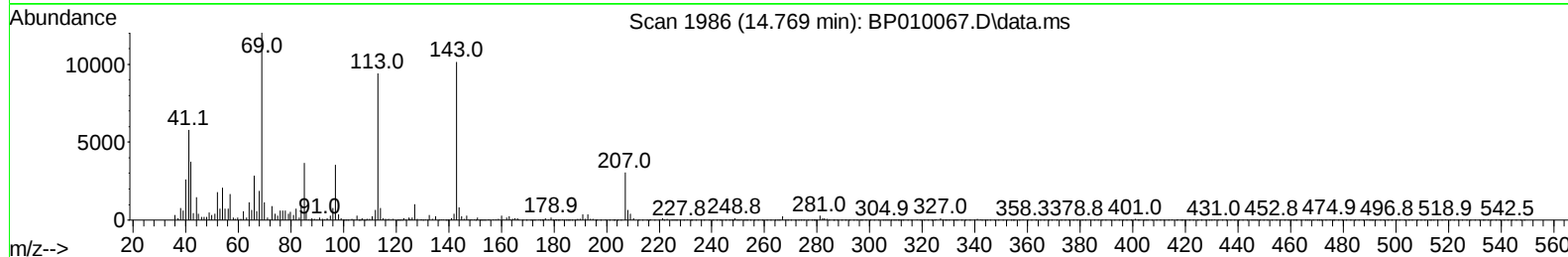
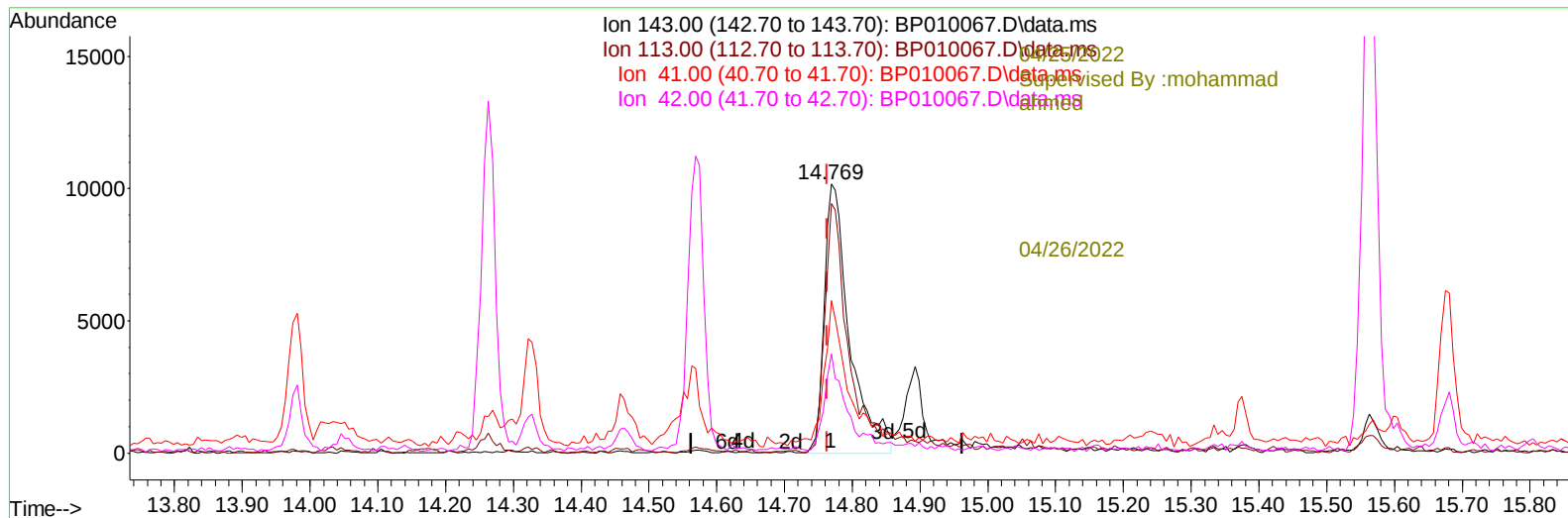
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ALS Vial : 63 Sample Multiplier: 1

Instrument :
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TIC: BP010067.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.769min (+ 0.006) 3.81 ng/ul m

response 26542

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	92.76#
41.00	23.20	56.79#
42.00	18.00	37.02#

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 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
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 C0J44

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/25/2022
 Supervised By :mohammad ahmed 04/26/2022

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----04/25/2022							
Supervised By :mohammad							
ahmed							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.993	152		143349	20.000	ng/ul	0.05
20) Naphthalene-d8	10.757	136		660027	20.000	ng/ul	0.01
38) Acenaphthene-d10	14.569	164		486521	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188		1068587	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240		784019	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264		660134	20.000	ng/ul	0.00
-----04/26/2022							
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.376	96		17546m	5.337	ng/uL	0.00
4) Pyridine-d5	3.817	84		46780	5.058	ng/ul	0.03
7) Phenol-d5	7.105	99		96116	7.371	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.281	67		295637	37.786	ng/ul	0.01
11) 2-Chlorophenol-d4	7.481	132		284377	29.449	ng/ul	0.01
15) 4-Methylphenol-d8	8.652	113		185634m	17.183	ng/ul	0.00
21) Nitrobenzene-d5	9.110	128		193363	36.835	ng/ul	0.01
24) 2-Nitrophenol-d4	9.828	143		182206	31.781	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.369	165		345075	30.918	ng/ul	0.00
31) 4-Chloroaniline-d4	10.881	131		303498	18.965	ng/ul	0.00
46) Dimethylphthalate-d6	13.981	166		1374732	36.324	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160		1556015	33.971	ng/ul	0.00
54) 4-Nitrophenol-d4	14.769	143		26542m	3.810	ng/ul	0.00
60) Fluorene-d10	15.563	176		1204461	36.997	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200		97190	14.384	ng/ul	0.00
73) Anthracene-d10	17.422	188		1895468	36.665	ng/ul	0.00
81) Pyrene-d10	19.651	212		2109609	47.314	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264		1311140	37.886	ng/ul	0.00
Target Compounds							
						Qvalue	
22) Nitrobenzene	9.163	77		4121892	320.480	ng/ul#	87
39) 1,2,4,5-Tetrachloroben...	12.769	216		626559	41.684	ng/ul	94
75) 1,2,3,4-Tetrachloroben...	13.375	216		818236	52.450	ng/uL#	87
76) Pentachlorobenzene	14.893	250		38643	2.526	ng/uL#	89

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

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