

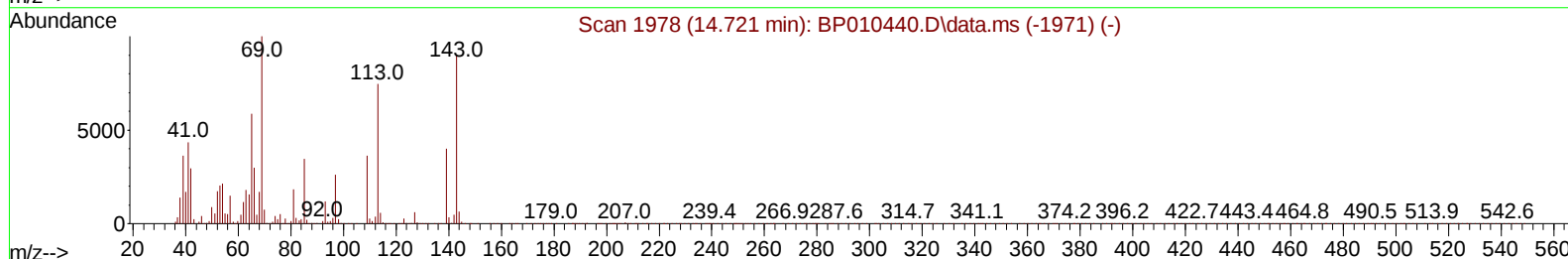
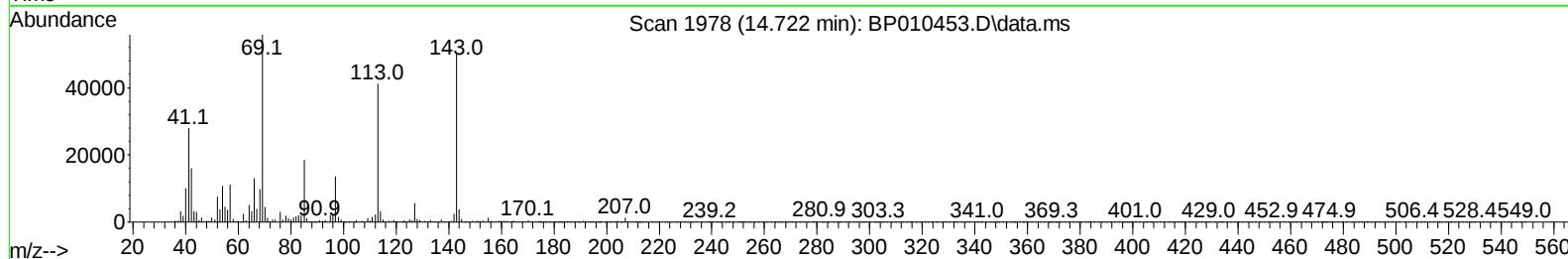
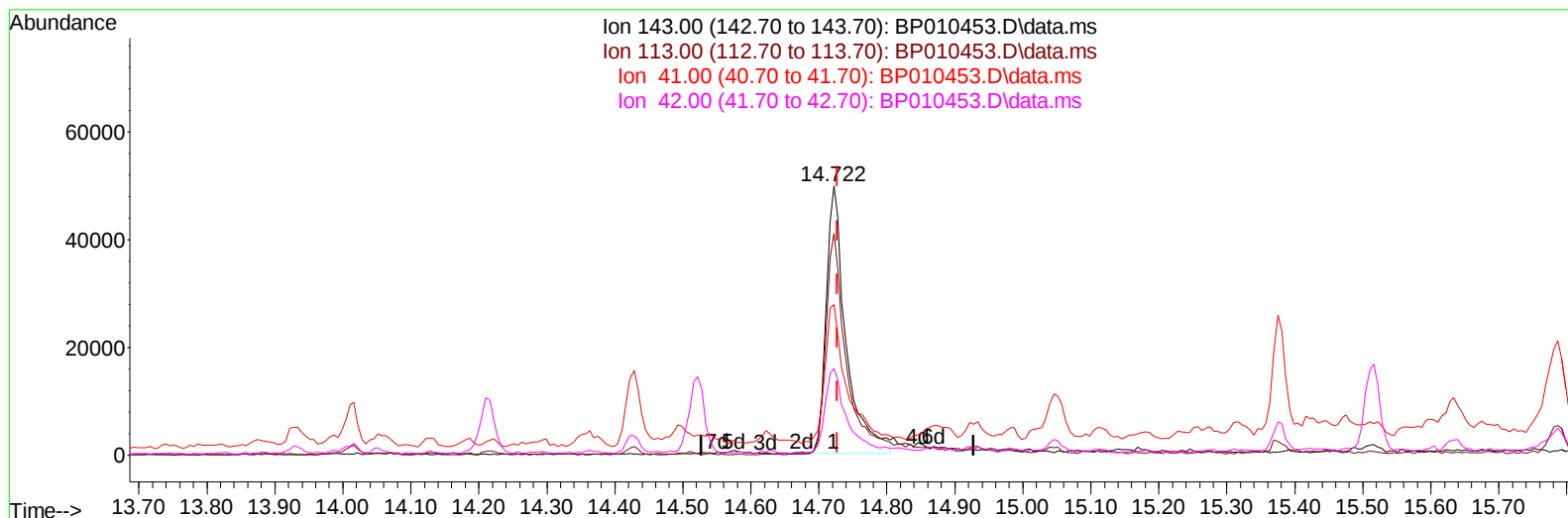
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010453.D
 Acq On : 21 May 2022 01:22
 Operator : CG/JU
 Sample : N2843-15
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW4J5

Manual IntegrationsAPPROVED

Quant Time: May 21 03:49:58 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 05/23/2022
 Supervised By :mohammad ahmed 05/25/2022



TIC: BP010453.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 12.56 ng/u1

response 101910

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.44#
41.00	23.20	55.93#
42.00	18.00	32.24#

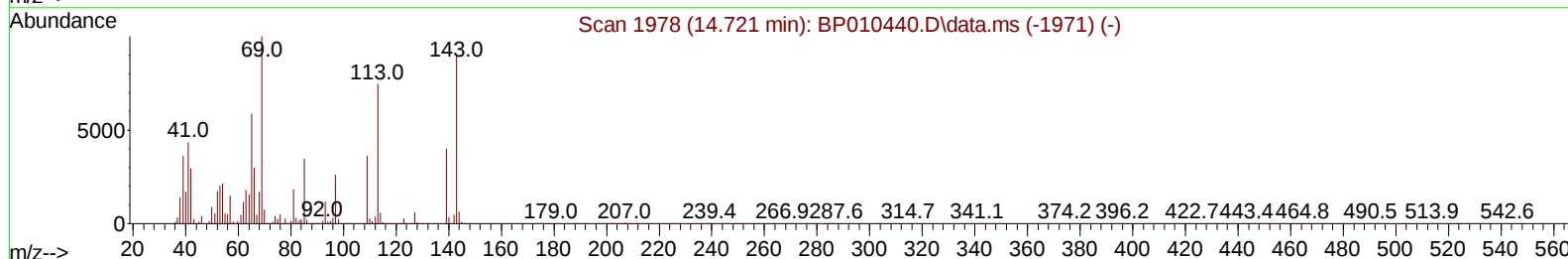
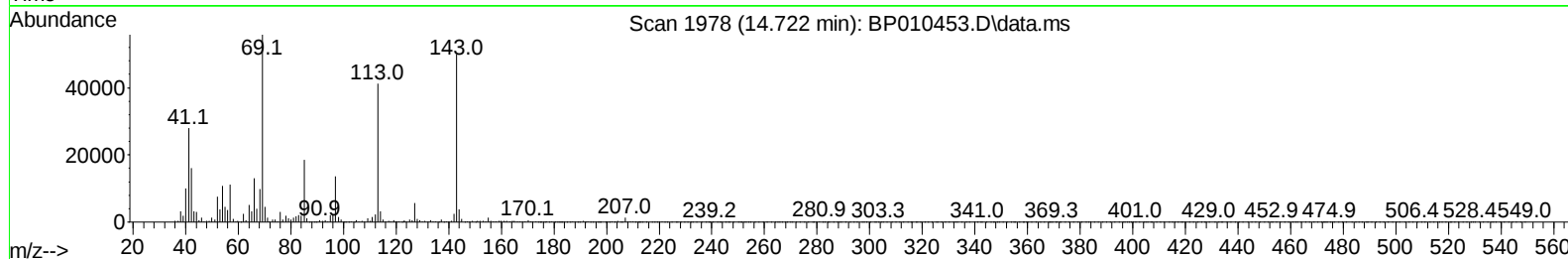
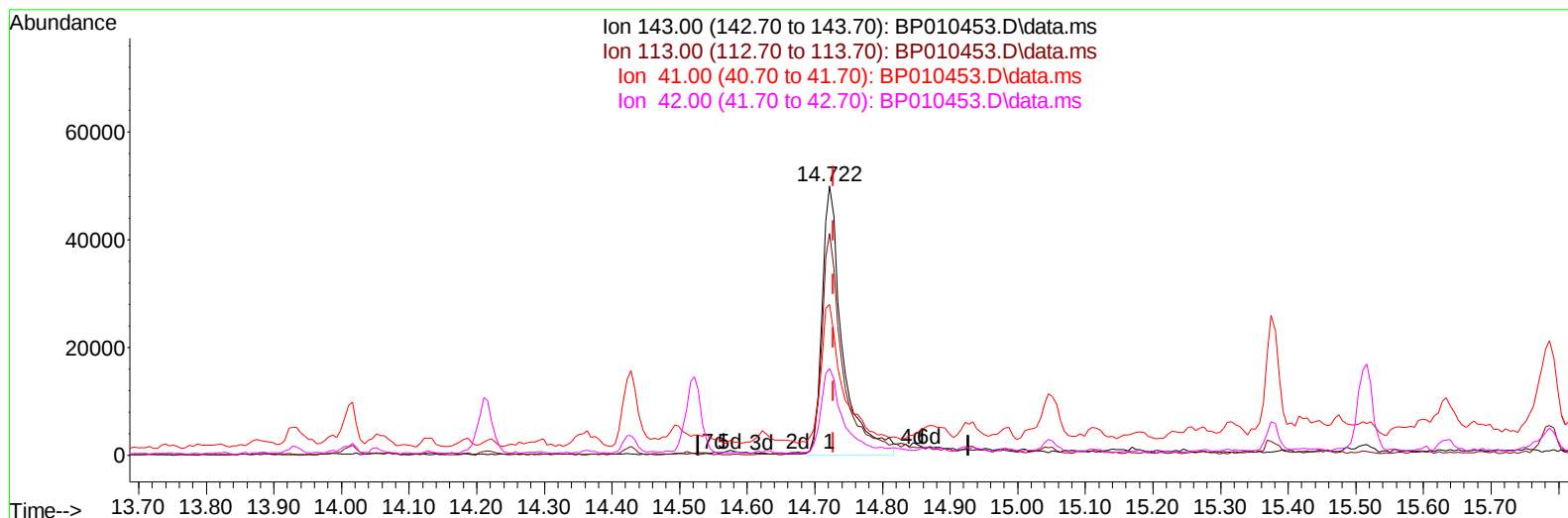
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TIC: BP010453.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 13.03 ng/ul m

response 105740

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.44#
41.00	23.20	55.93#
42.00	18.00	32.24#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.887	152	199957	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.687	136	903294	20.000	ng/ul	#-0.02
38) Acenaphthene-d10	14.522	164	618924	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	1272235	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	1144025	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	1053636	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	17654	3.763	ng/uL	0.00
4) Pyridine-d5	3.764	84	120943	9.080	ng/ul	0.00
7) Phenol-d5	7.058	99	337929	20.135	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.216	67	238101	21.116	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.422	132	280696	21.857	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	287062	20.127	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	148141	21.191	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.775	143	157897	21.312	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	277344	19.798	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	396588	19.040	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	968991	21.250	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1139083	21.656	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	105740m	13.028	ng/ul	0.00
60) Fluorene-d10	15.516	176	858346	21.636	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.634	200	91043	12.236	ng/ul	-0.01
73) Anthracene-d10	17.375	188	1316912	22.860	ng/ul	0.00
81) Pyrene-d10	19.604	212	1529662	25.177	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.627	264	1243212	23.498	ng/ul	0.00
Target Compounds						
					Qvalue	
16) Acetophenone	8.716	105	23832	1.063	ng/ul#	67

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

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