

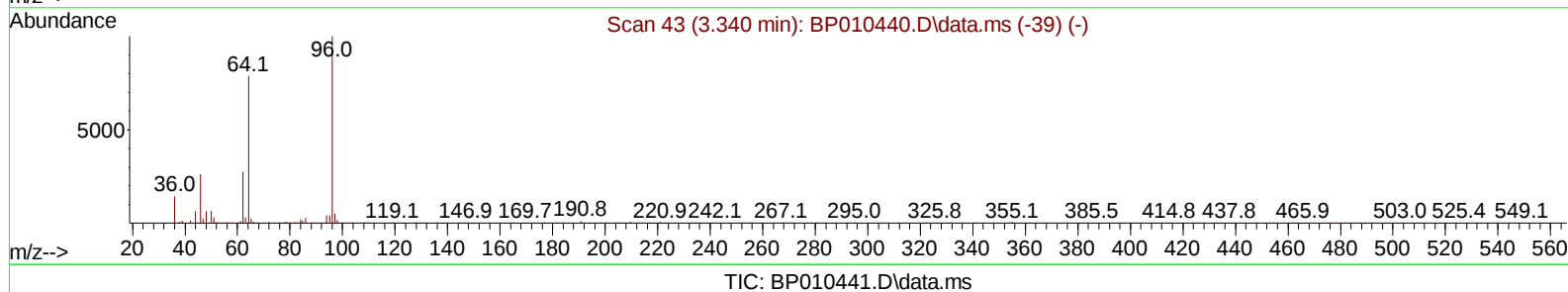
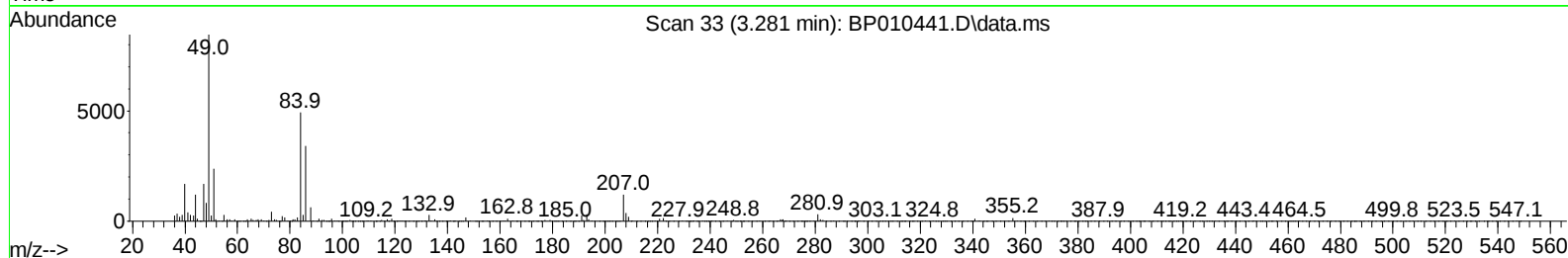
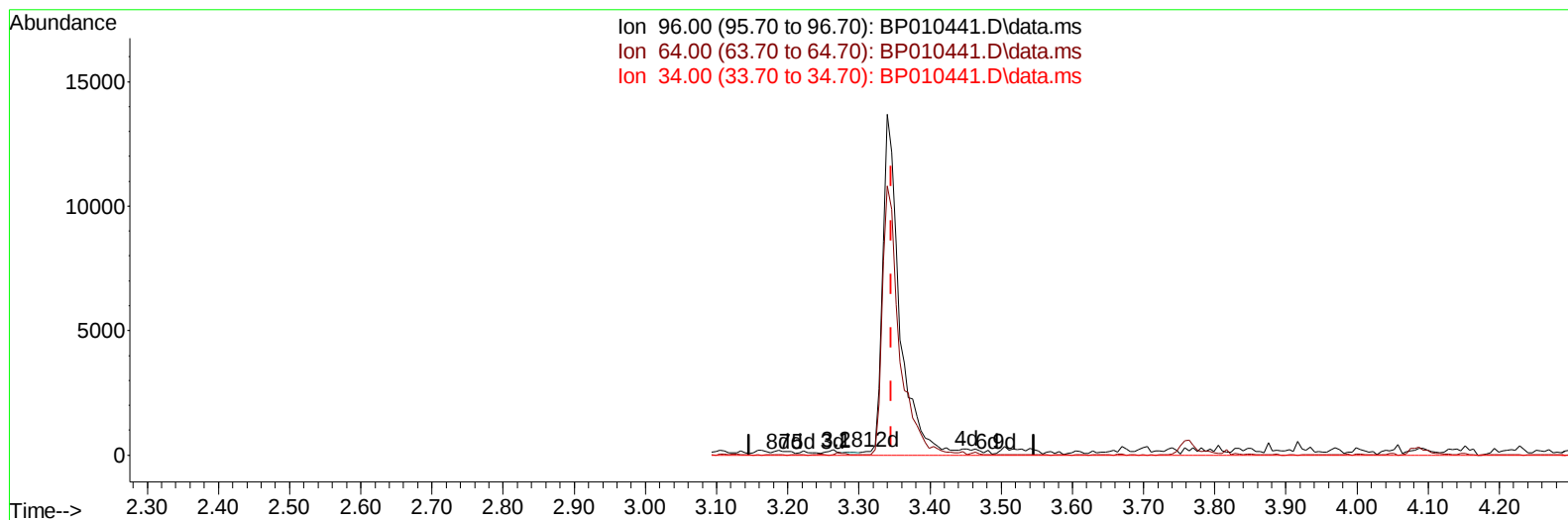
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052022\
 Data File : BP010441.D
 Acq On : 20 May 2022 18:36
 Operator : CG/JU
 Sample : PB144914BL
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK914

Manual IntegrationsAPPROVED

Quant Time: May 20 23:03:28 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



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

3.281min (-0.065) 0.01 ng/uL

response 22

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	55.40	60.94
34.00	0.00	0.00
0.00	0.00	0.00

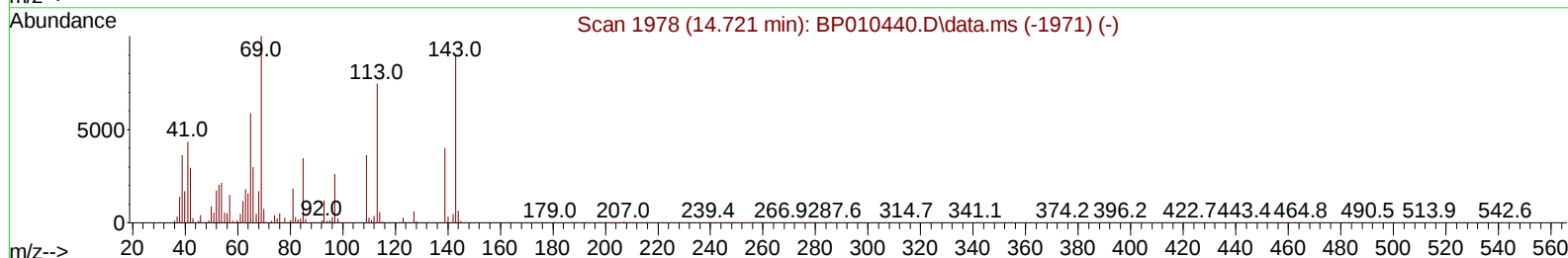
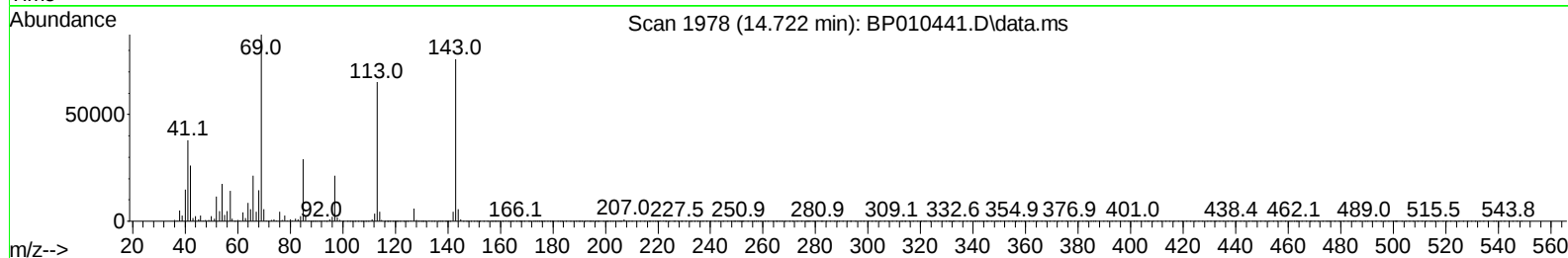
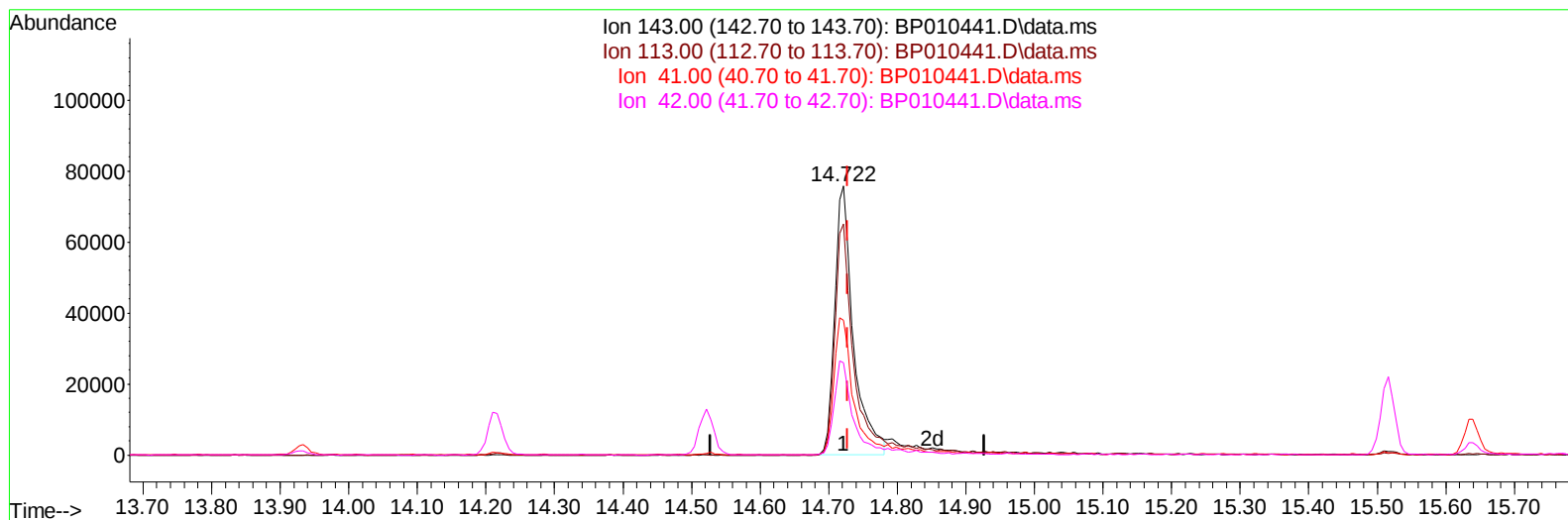
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TIC: BP010441.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 22.13 ng/u1

response 142428

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.94#
41.00	23.20	50.05#
42.00	18.00	34.30#

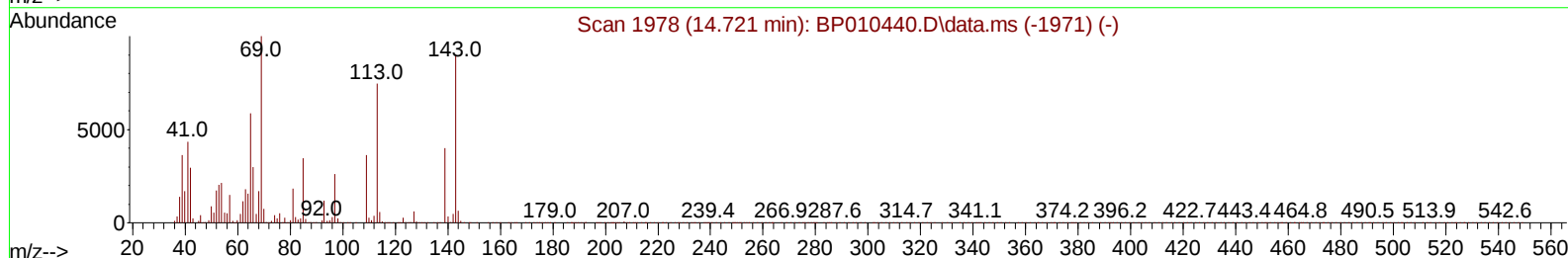
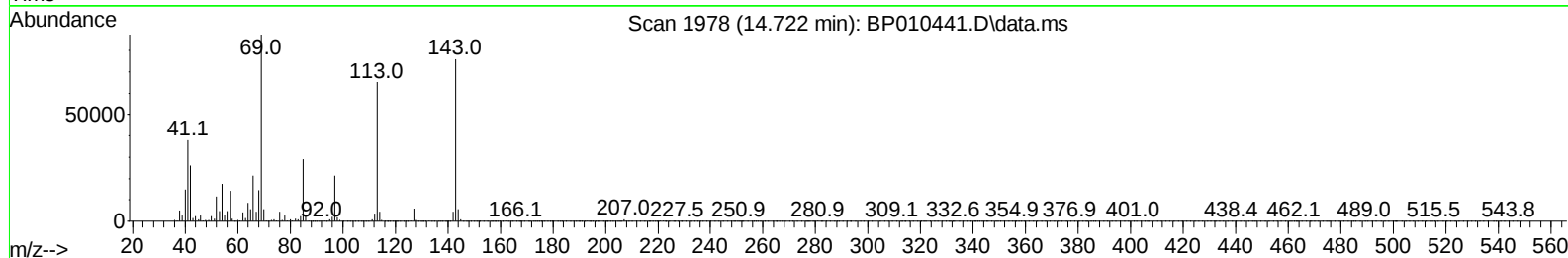
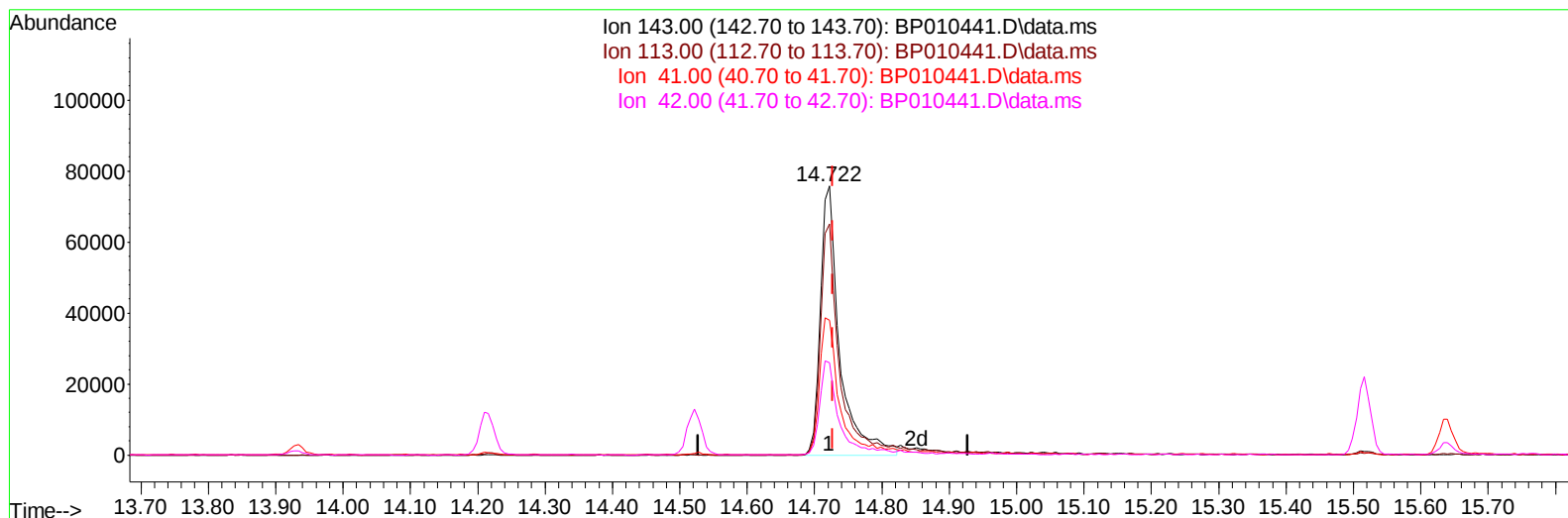
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Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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TIC: BP010441.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.722min (-0.006) 23.46 ng/ul m

response 150937

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	85.94#
41.00	23.20	50.05#
42.00	18.00	34.30#

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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	157195	20.000	ng/ul	0.00
20) Naphthalene-d8	10.693	136	707211	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.522	164	490710	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.275	188	1057007	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.357	240	964169	20.000	ng/ul	# 0.00
88) Perylene-d12	23.780	264	804142	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.340	96	21890m	5.936	ng/uL	0.00
4) Pyridine-d5	3.758	84	317632	30.333	ng/ul	0.00
7) Phenol-d5	7.058	99	393101	29.795	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.222	67	278580	31.427	ng/ul	0.00
11) 2-Chlorophenol-d4	7.422	132	314893	31.191	ng/ul	-0.01
15) 4-Methylphenol-d8	8.599	113	339665	30.293	ng/ul	-0.01
21) Nitrobenzene-d5	9.052	128	171765	31.382	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	185509	31.982	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.310	165	320447	29.217	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.828	131	474032	29.068	ng/ul	-0.01
46) Dimethylphthalate-d6	13.934	166	1163896	32.193	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1301462	31.208	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	150937m	23.455	ng/ul	0.00
60) Fluorene-d10	15.516	176	980911	31.186	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.640	200	149007	24.104	ng/ul	0.00
73) Anthracene-d10	17.375	188	1571637	32.836	ng/ul	0.00
81) Pyrene-d10	19.604	212	1855085	36.229	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.627	264	1347767	33.378	ng/ul	0.00

Target Compounds Qvalue

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

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