

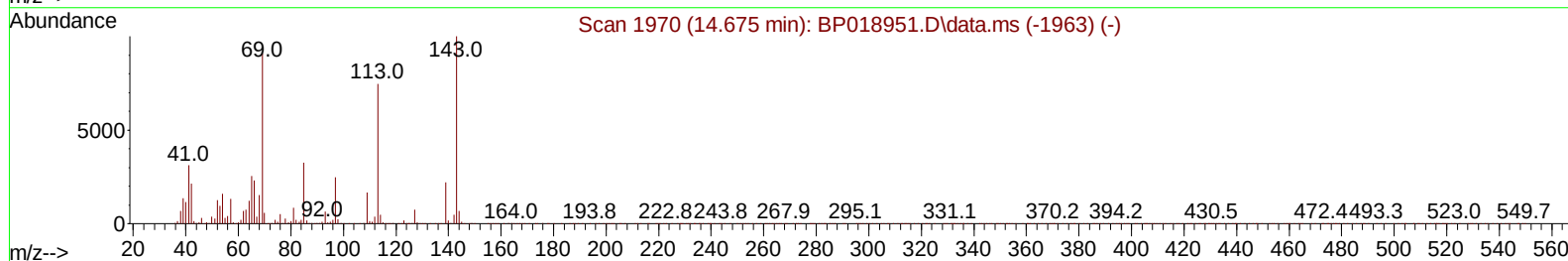
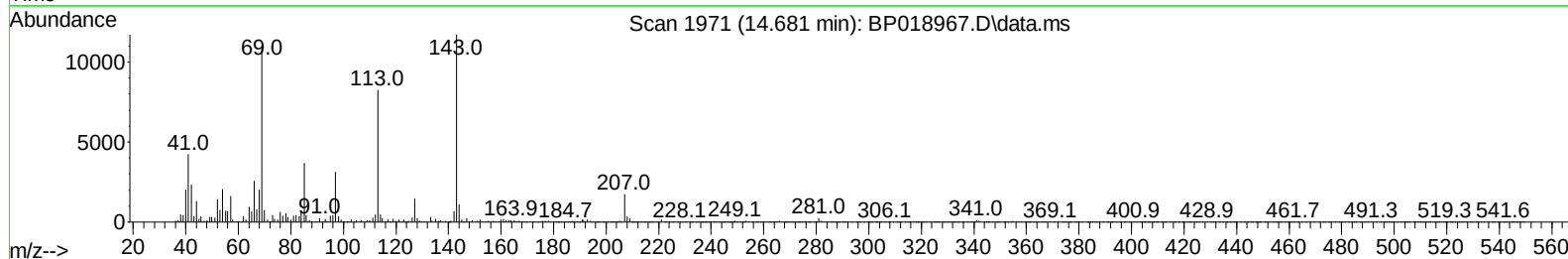
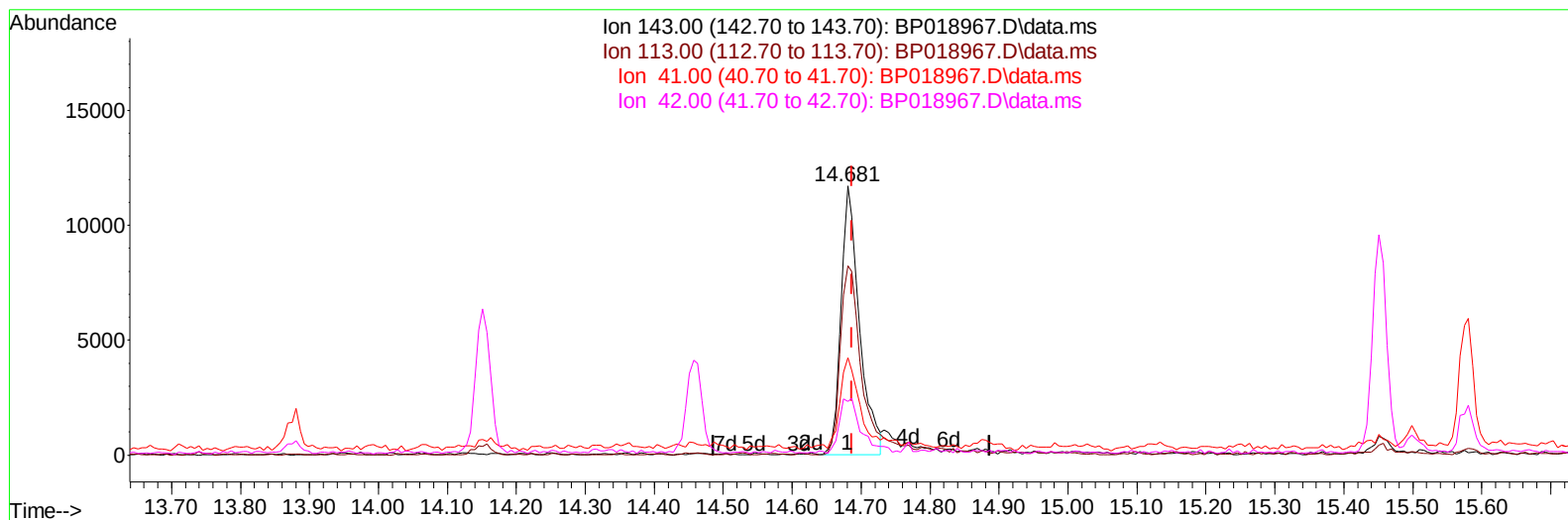
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122023\
 Data File : BP018967.D
 Acq On : 20 Dec 2023 18:53
 Operator : MA/JU
 Sample : 05894-05
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH3K4

Manual IntegrationsAPPROVED

Quant Time: Dec 20 21:50:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 19 05:28:03 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/21/2023
 Supervised By :mohammad ahmed 12/22/2023



TIC: BP018967.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (-0.006) 5.65 ng/ul

response 21634

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	70.48
41.00	34.60	36.25
42.00	23.00	20.03

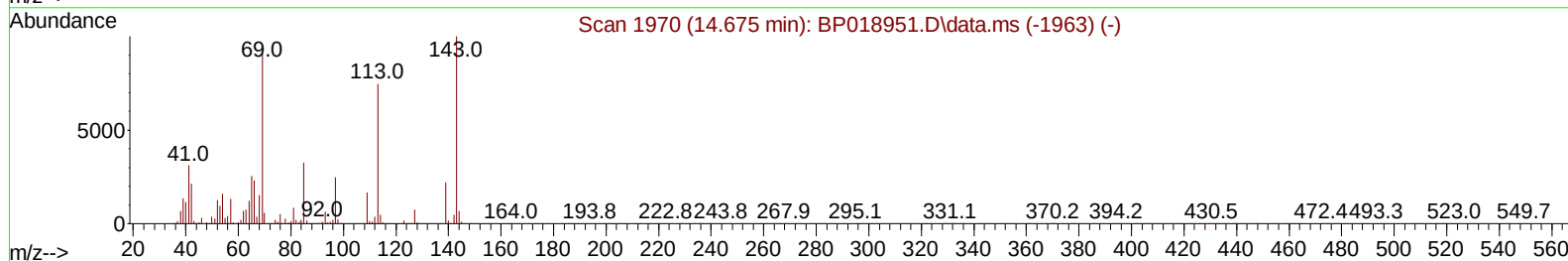
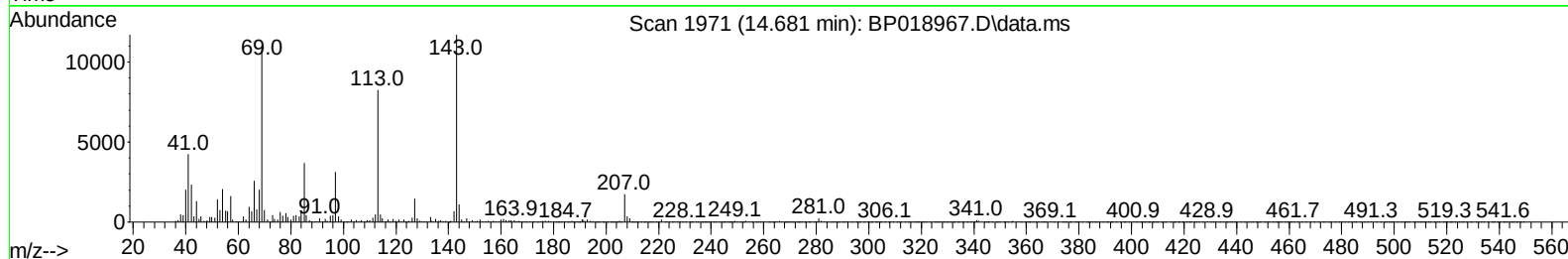
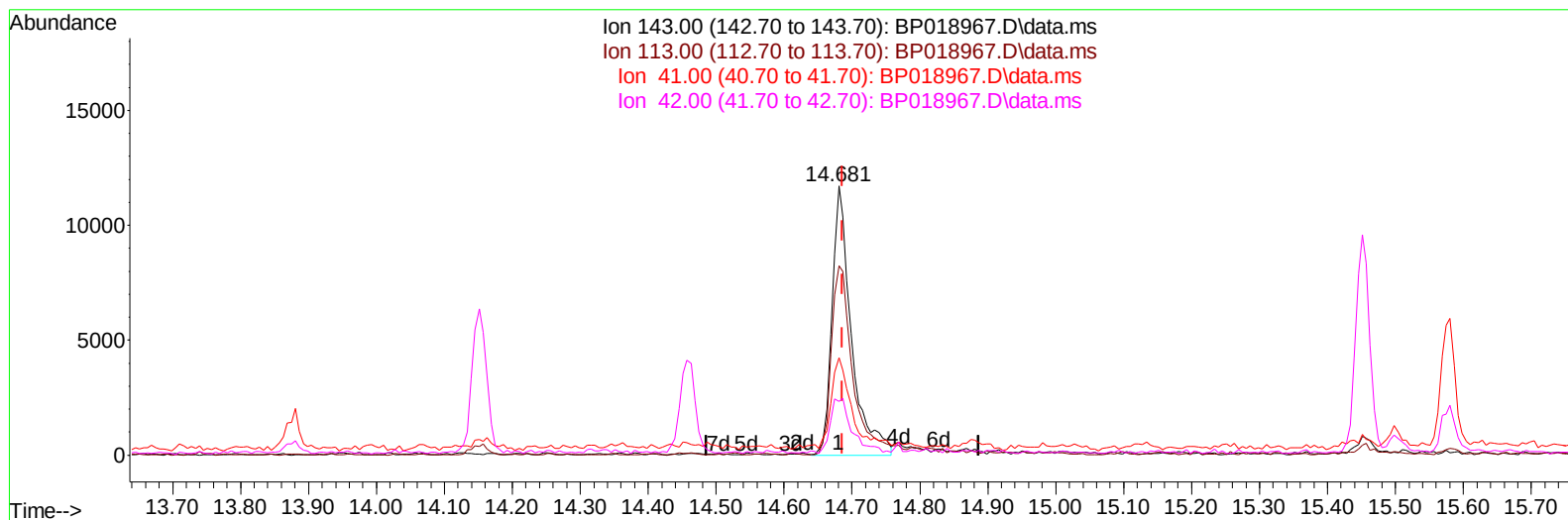
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TIC: BP018967.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (-0.006) 6.04 ng/ul m

response 23103

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	70.48
41.00	34.60	36.25
42.00	23.00	20.03

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/21/2023
 Supervised By :mohammad ahmed 12/22/2023

Quant Time: Dec 20 22:45:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	128870	20.000	ng/ul	0.00
20) Naphthalene-d8	10.622	136	445617	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	261273	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.210	188	602427	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.298	240	743072	20.000	ng/ul	-0.02
88) Perylene-d12	23.662	264	928438	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	15476	4.092	ng/uL	0.00
4) Pyridine-d5	3.675	84	102228	10.092	ng/ul	0.00
7) Phenol-d5	6.999	99	64835	5.013	ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth...	7.163	67	205940	26.868	ng/ul	0.00
11) 2-Chlorophenol-d4	7.357	132	211032	20.912	ng/ul	-0.01
15) 4-Methylphenol-d8	8.540	113	125376	11.979	ng/ul	-0.02
21) Nitrobenzene-d5	8.987	128	132223	33.338	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.704	143	140806	34.243	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.246	165	242046	28.787	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.763	131	221204	20.131	ng/ul	-0.01
46) Dimethylphthalate-d6	13.875	166	842014	36.109	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	920754	35.859	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143	23103m	6.036	ng/ul	0.00
60) Fluorene-d10	15.451	176	732239	37.409	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.581	200	137777	38.142	ng/ul	-0.01
73) Anthracene-d10	17.310	188	1262452	39.862	ng/ul	0.00
81) Pyrene-d10	19.545	212	1683799	29.110	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.510	264	2123173	39.220	ng/ul	-0.02

Target Compounds Qvalue

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

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