

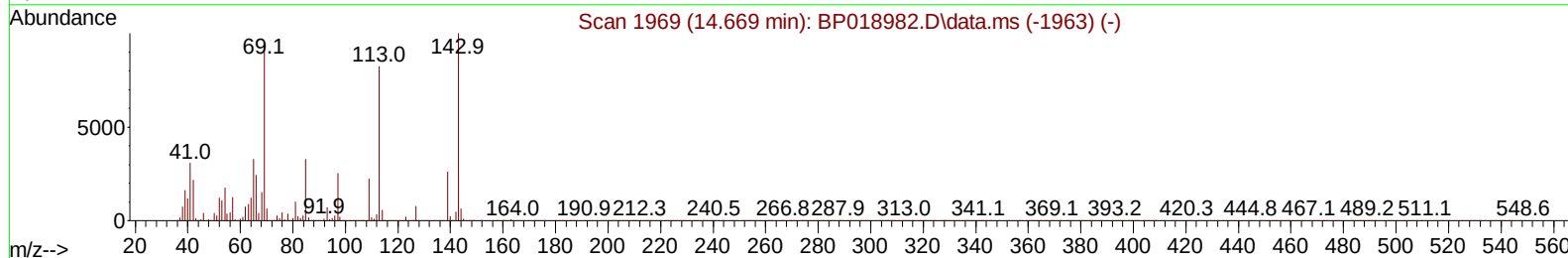
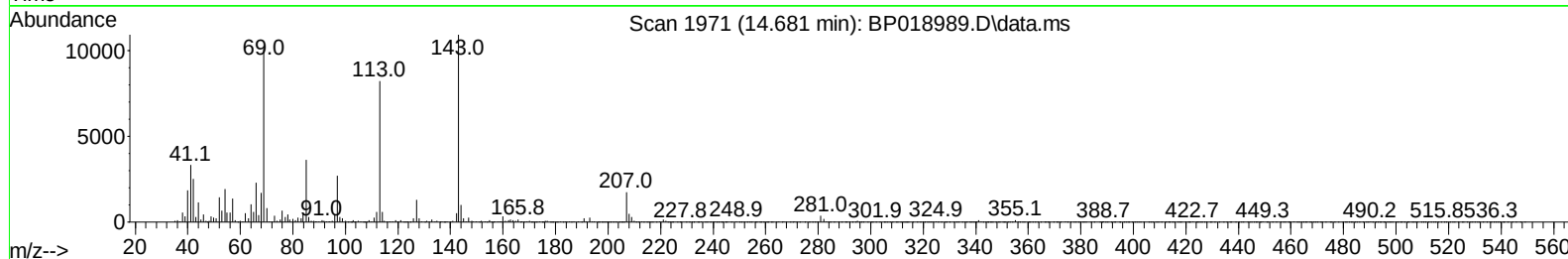
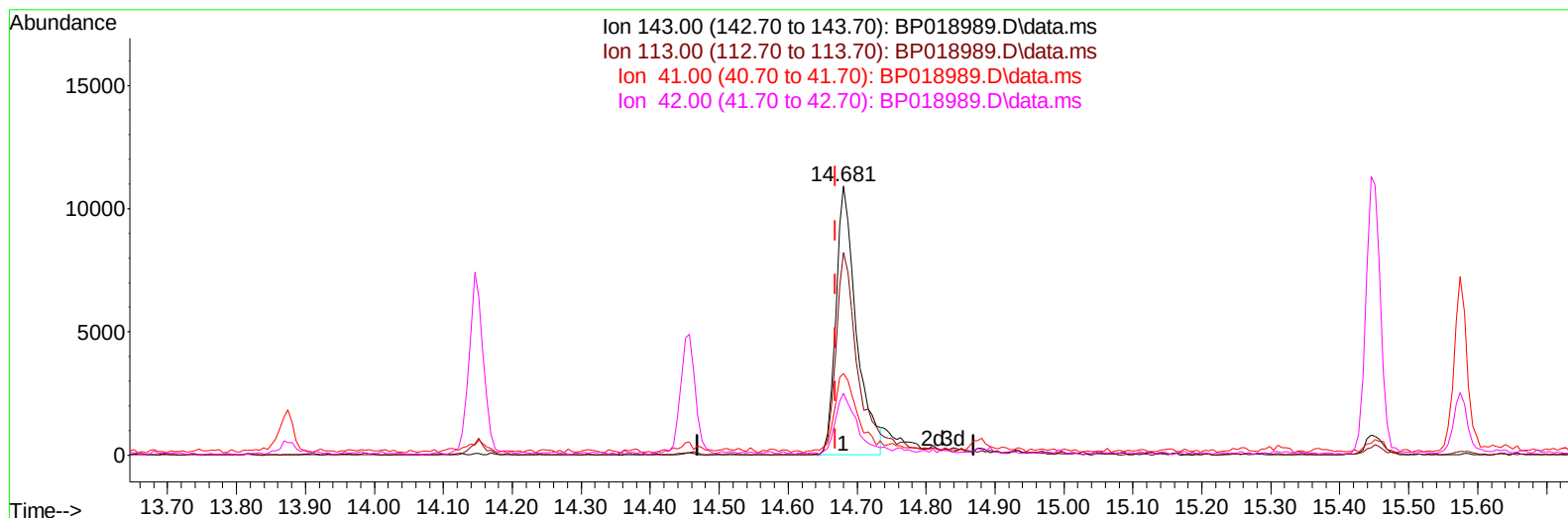
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122023\
Data File : BP018989.D
Acq On : 21 Dec 2023 07:52
Operator : MA/JU
Sample : 05894-14
Misc :
ALS Vial : 39 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH3M0

Manual IntegrationsAPPROVED

Quant Time: Dec 21 08:19:56 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Dec 21 04:46:12 2023
Response via : Initial Calibration

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



TIC: BP018989.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (+ 0.012) 5.50 ng/ul

response 22434

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	75.36
41.00	34.60	30.40
42.00	23.00	23.10

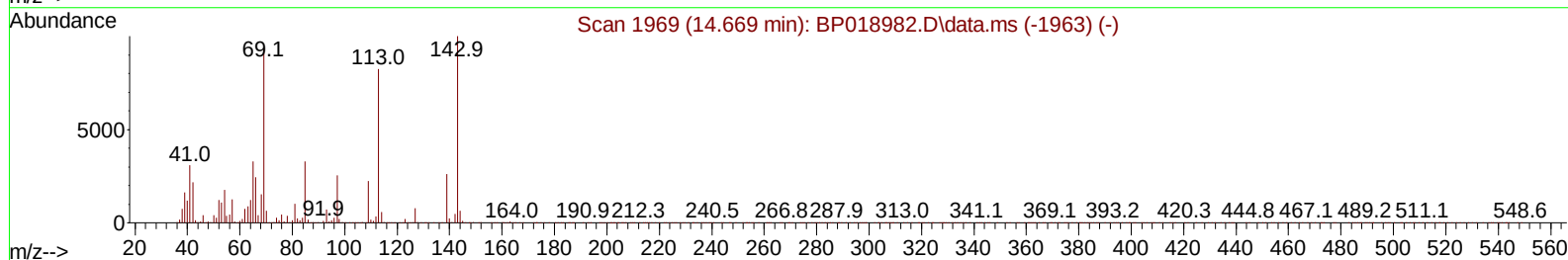
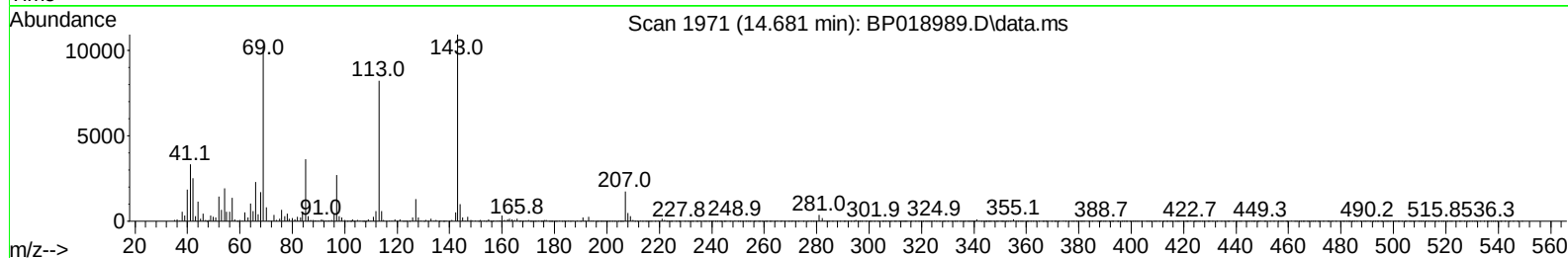
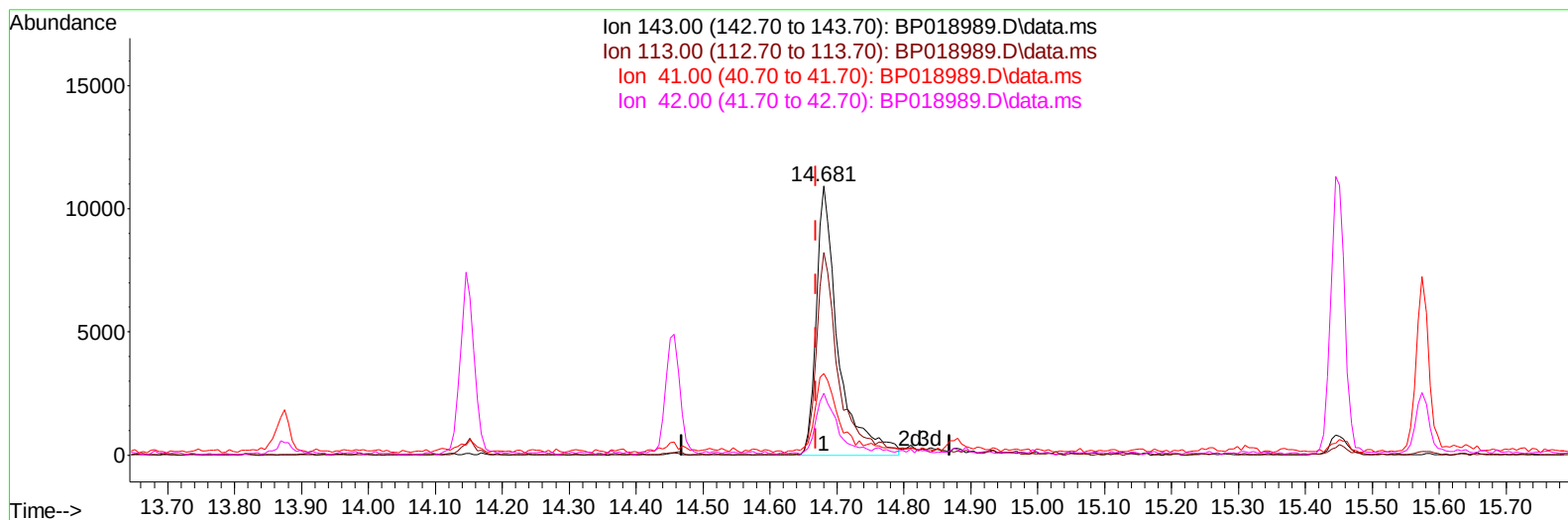
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TIC: BP018989.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (+ 0.012) 6.07 ng/ul m

response 24766

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	75.36
41.00	34.60	30.40
42.00	23.00	23.10

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Manual IntegrationsAPPROVED

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

Quant Time: Dec 21 08:21:02 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	110770	20.000	ng/ul	0.00
20) Naphthalene-d8	10.616	136	417864	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	278508	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	683941	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	841403	20.000	ng/ul	# 0.00
88) Perylene-d12	23.657	264	1011986	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	13070	4.021	ng/uL	0.00
4) Pyridine-d5	3.669	84	93401	10.727	ng/ul	0.00
7) Phenol-d5	6.999	99	59045	5.312	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.157	67	196100	29.764	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	192923	22.242	ng/ul	0.00
15) 4-Methylphenol-d8	8.534	113	118776	13.203	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	129006	34.687	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	140123	36.340	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	244675	31.033	ng/ul	0.00
31) 4-Chloroaniline-d4	10.757	131	320708	31.125	ng/ul	0.00
46) Dimethylphthalate-d6	13.875	166	960175	38.628	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	1018277	37.203	ng/ul	0.00
54) 4-Nitrophenol-d4	14.681	143	24766m	6.070	ng/ul	0.01
60) Fluorene-d10	15.451	176	856972	41.072	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	159150	38.808	ng/ul	0.00
73) Anthracene-d10	17.304	188	1521878	42.326	ng/ul	0.00
81) Pyrene-d10	19.545	212	2040379	31.153	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.510	264	2527428	42.833	ng/ul	0.00

Target Compounds Qvalue

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

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