

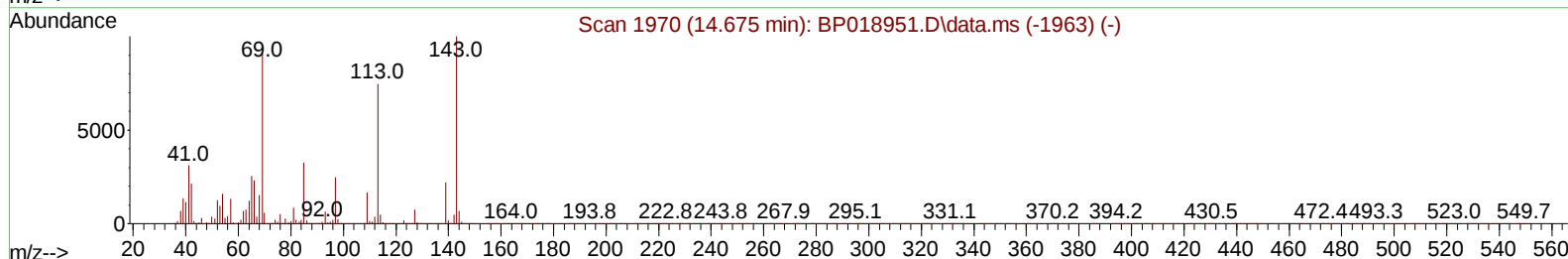
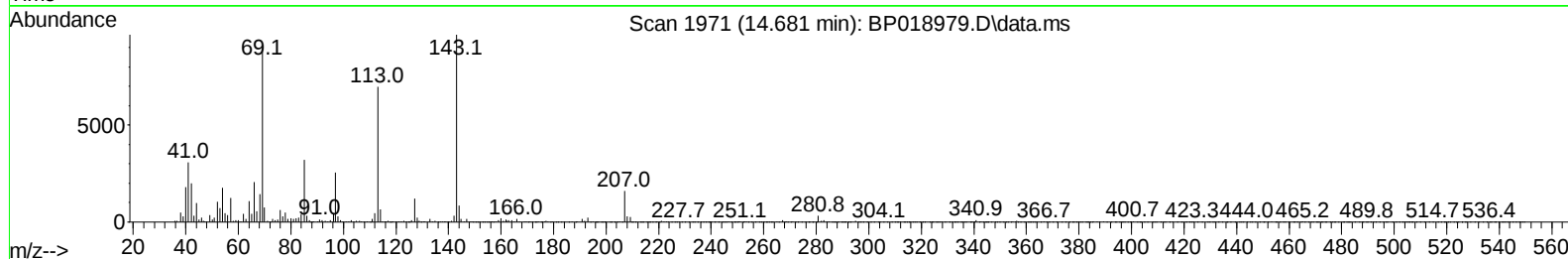
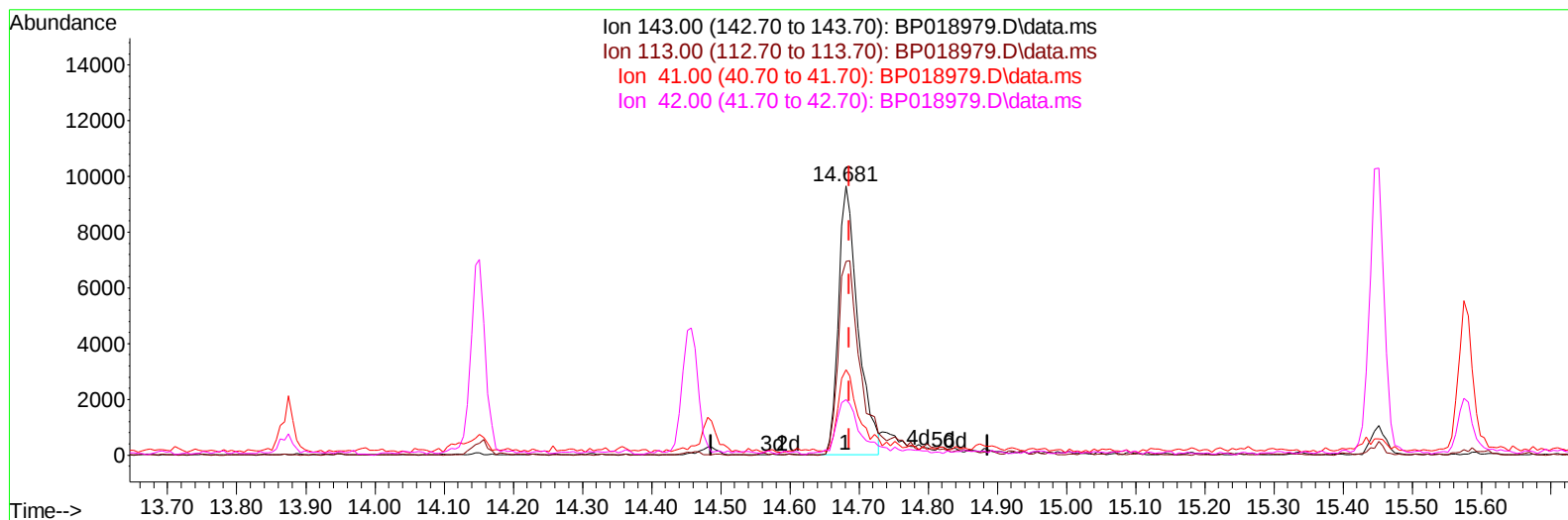
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122023\
 Data File : BP018979.D
 Acq On : 21 Dec 2023 01:40
 Operator : MA/JU
 Sample : 05894-12
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH3L8

Manual IntegrationsAPPROVED

Quant Time: Dec 21 03:19:39 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: BP018979.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (-0.006) 4.33 ng/u1

response 18874

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	72.06
41.00	34.60	31.79
42.00	23.00	20.66

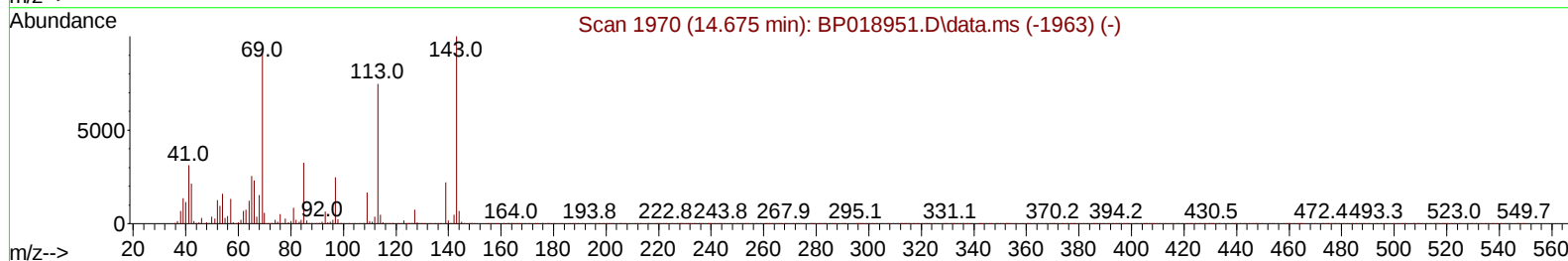
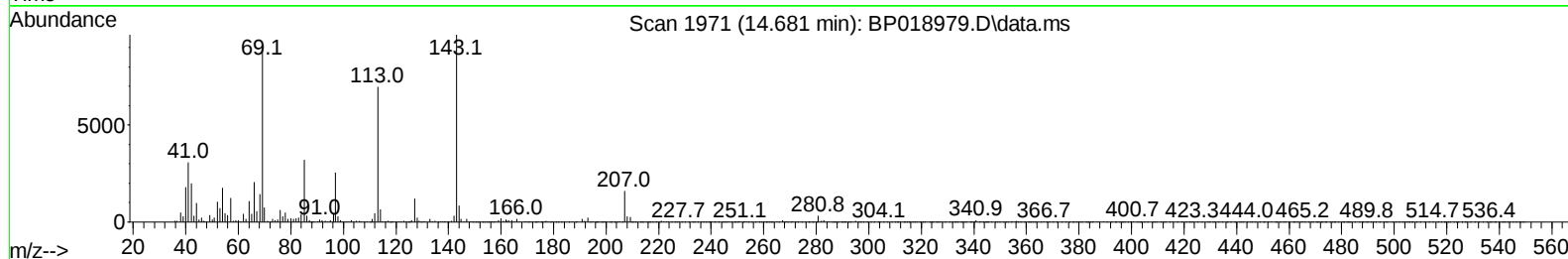
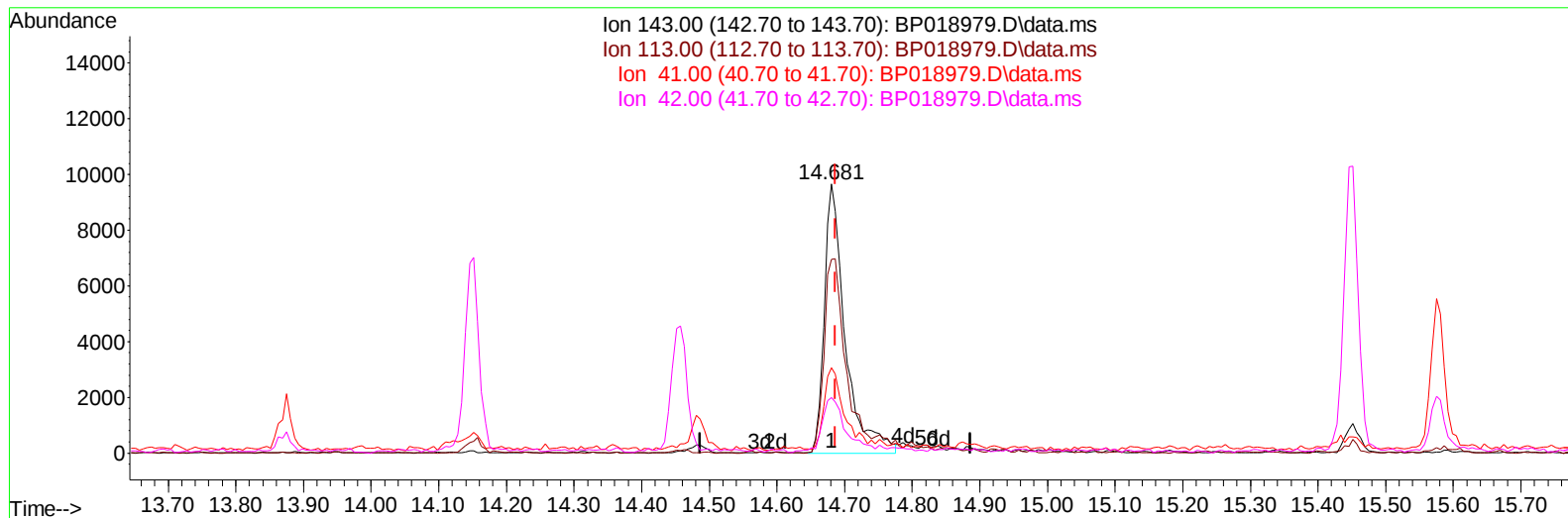
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TIC: BP018979.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.681min (-0.006) 4.73 ng/ul m

response 20612

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	70.80	72.06
41.00	34.60	31.79
42.00	23.00	20.66

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/21/2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.822	152	117085	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.616	136	447777	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.457	164	297307	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.210	188	711314	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.298	240	828945	20.000	ng/ul	#-0.02
88) Perylene-d12	23.663	264	1001420	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	11490	3.344	ng/uL	0.00
4) Pyridine-d5	3.675	84	93141	10.120	ng/ul	0.00
7) Phenol-d5	6.999	99	59300	5.047	ng/ul	-0.02
9) Bis-(2-Chloroethyl)eth...	7.163	67	191655	27.521	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	192234	20.967	ng/ul	-0.02
15) 4-Methylphenol-d8	8.540	113	117333	12.339	ng/ul	-0.02
21) Nitrobenzene-d5	8.987	128	128960	32.358	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.705	143	136705	33.085	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.240	165	243322	28.799	ng/ul	-0.02
31) 4-Chloroaniline-d4	10.763	131	319224	28.911	ng/ul	-0.01
46) Dimethylphthalate-d6	13.875	166	890004	33.541	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	955205	32.692	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.681	143	20612m	4.732	ng/ul	0.00
60) Fluorene-d10	15.451	176	775089	34.799	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	139257	32.651	ng/ul	-0.02
73) Anthracene-d10	17.310	188	1314857	35.161	ng/ul	0.00
81) Pyrene-d10	19.545	212	1730757	26.822	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.510	264	2043455	34.997	ng/ul	-0.02

Target Compounds Qvalue

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

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