

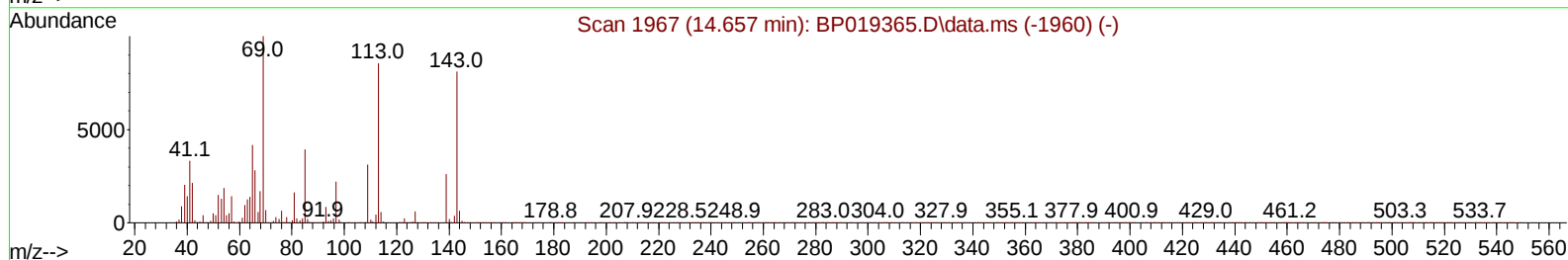
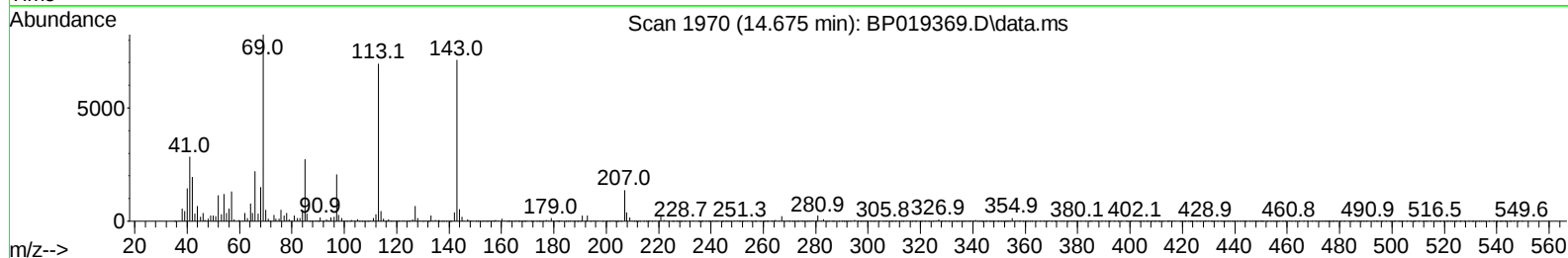
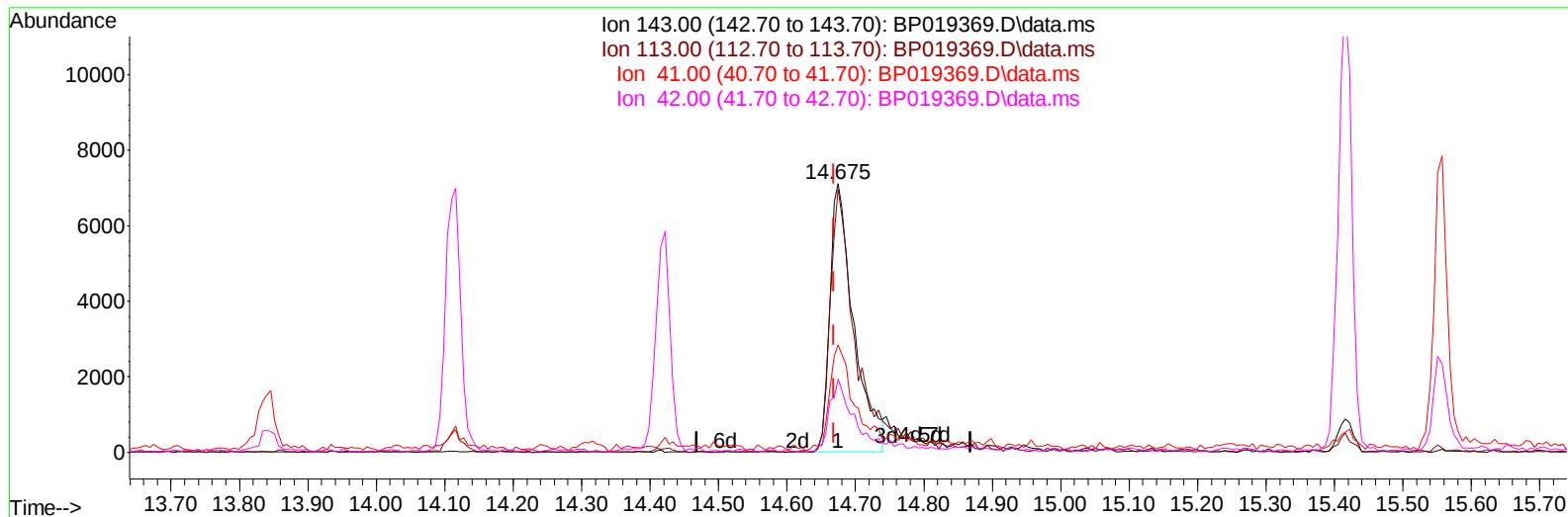
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011524\
Data File : BP019369.D
Acq On : 15 Jan 2024 12:25
Operator : MA/JU
Sample : P1130-06
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AZ4

Manual IntegrationsAPPROVED

Quant Time: Jan 15 12:51:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jan 09 03:05:24 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/16/2024
Supervised By :mohammad ahmed 01/17/2024



TIC: BP019369.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.006) 4.55 ng/u1

response 17071

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	97.92
41.00	39.60	39.96
42.00	25.70	27.36

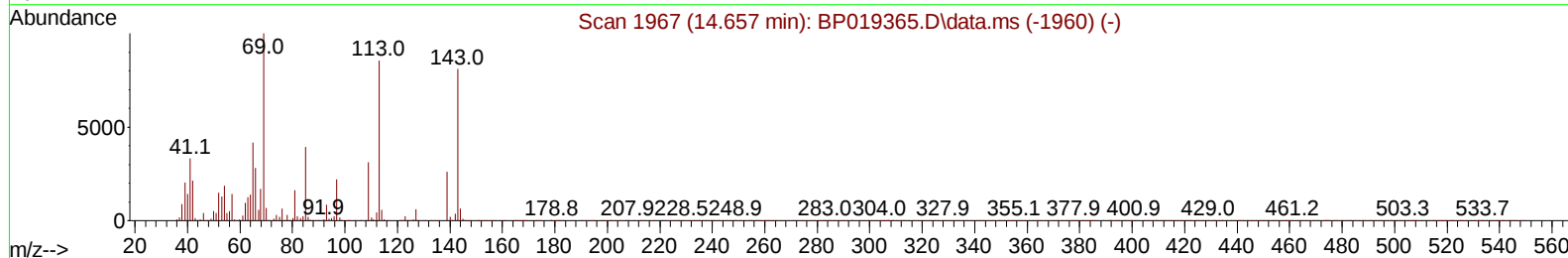
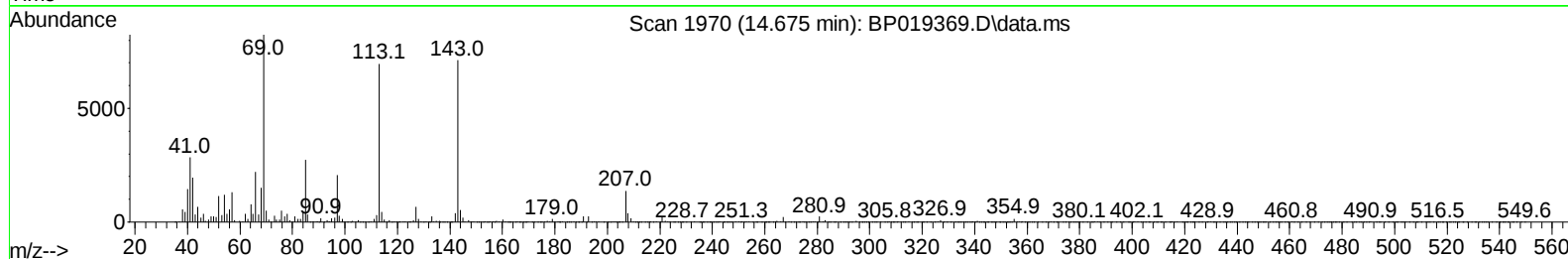
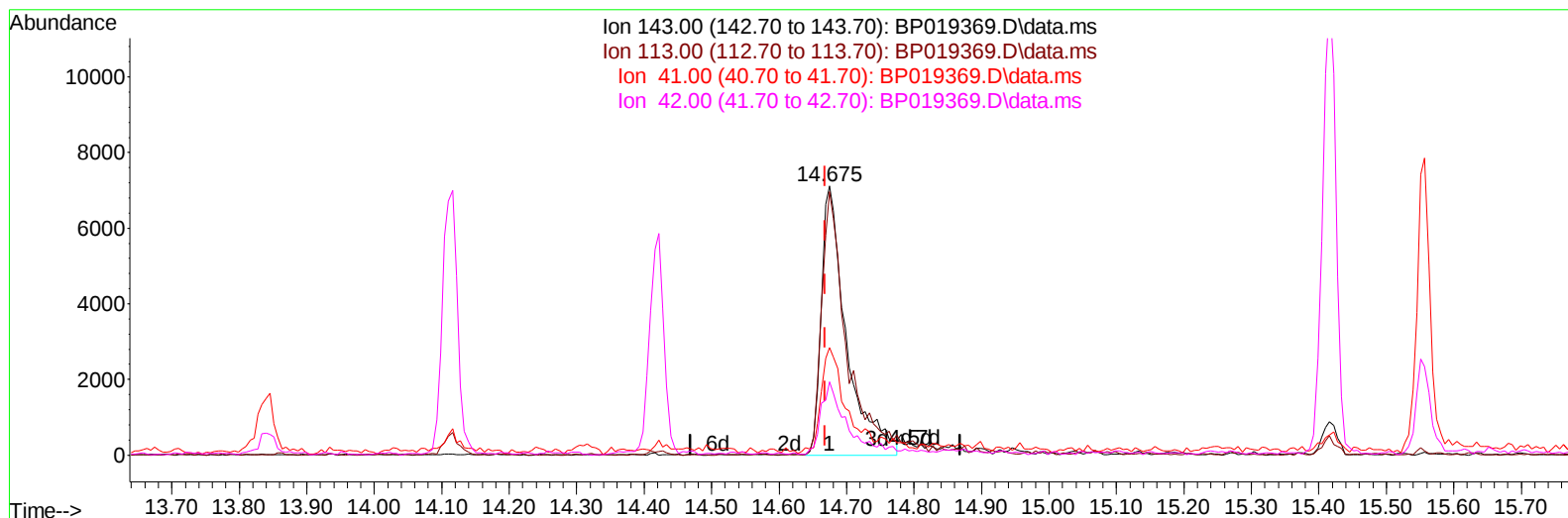
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TIC: BP019369.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.675min (+ 0.006) 4.89 ng/ul m

response 18335

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	97.92
41.00	39.60	39.96
42.00	25.70	27.36

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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 01/17/2024

Quant Time: Jan 16 00:26:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP010824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jan 09 03:05:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	98277	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.569	136	399061	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.422	164	276436	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.175	188	679525	20.000	ng/ul	-0.01
79) Chrysene-d12	21.280	240	653535	20.000	ng/ul	0.00
88) Perylene-d12	23.633	264	637147	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	11865	4.565	ng/uL	0.00
4) Pyridine-d5	3.640	84	83378	11.094	ng/ul	0.00
7) Phenol-d5	6.963	99	57766	6.150	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.116	67	191327	34.054	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.310	132	157112	24.349	ng/ul	0.00
15) 4-Methylphenol-d8	8.504	113	115540	15.505	ng/ul	-0.01
21) Nitrobenzene-d5	8.946	128	115610	34.436	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	128805	33.019	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.204	165	232153	29.457	ng/ul	0.00
31) 4-Chloroaniline-d4	10.722	131	320462	32.191	ng/ul	0.00
46) Dimethylphthalate-d6	13.839	166	853295	34.939	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	891557	34.055	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.675	143	18335m	4.889	ng/ul	0.00
60) Fluorene-d10	15.416	176	739828	37.001	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	159043	31.702	ng/ul	0.00
73) Anthracene-d10	17.275	188	1245714	36.757	ng/ul	-0.01
81) Pyrene-d10	19.522	212	1623376	38.094	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.480	264	1342831	38.760	ng/ul	0.00

Target Compounds Qvalue

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

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