

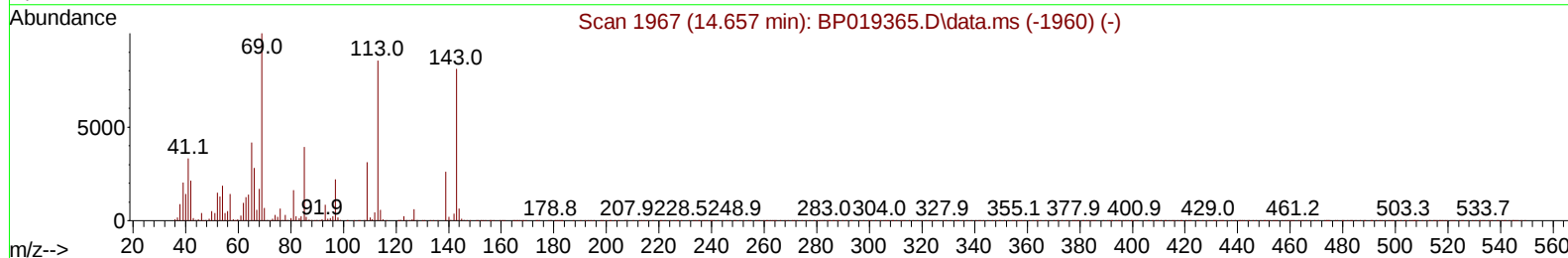
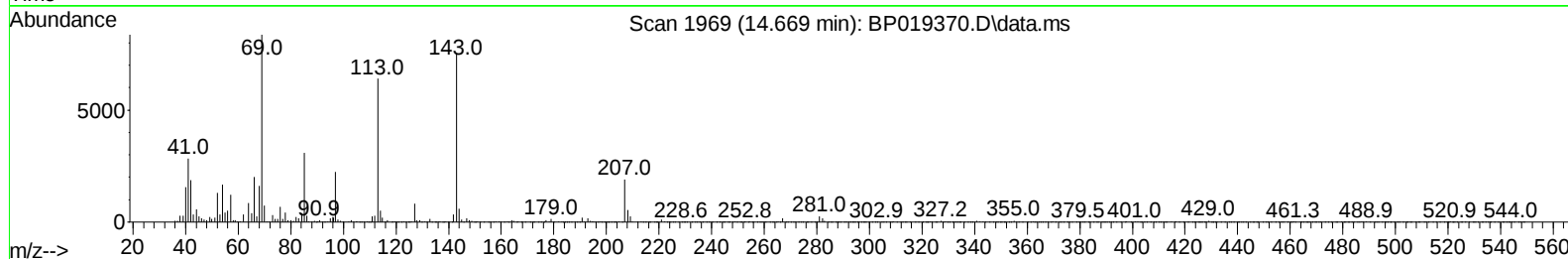
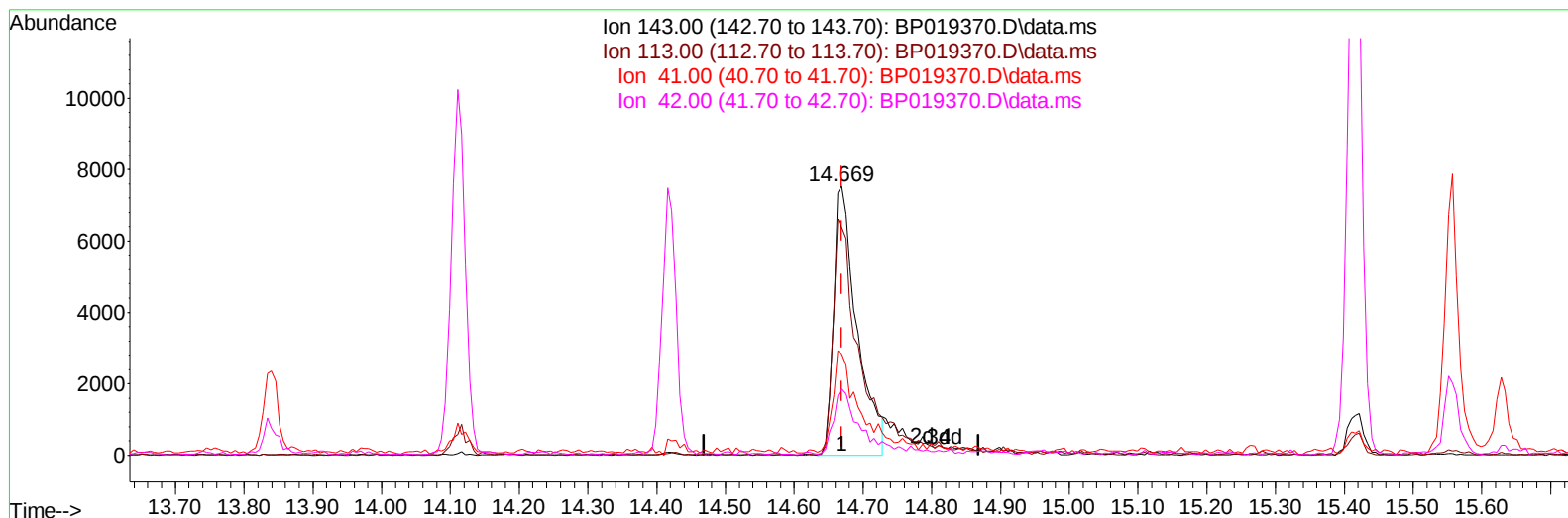
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011524\
Data File : BP019370.D
Acq On : 15 Jan 2024 12:59
Operator : MA/JU
Sample : P1130-07
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AZ8

Manual IntegrationsAPPROVED

Quant Time: Jan 15 14:11:15 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: BP019370.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 3.90 ng/u1

response 17751

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	84.85
41.00	39.60	37.71
42.00	25.70	24.66

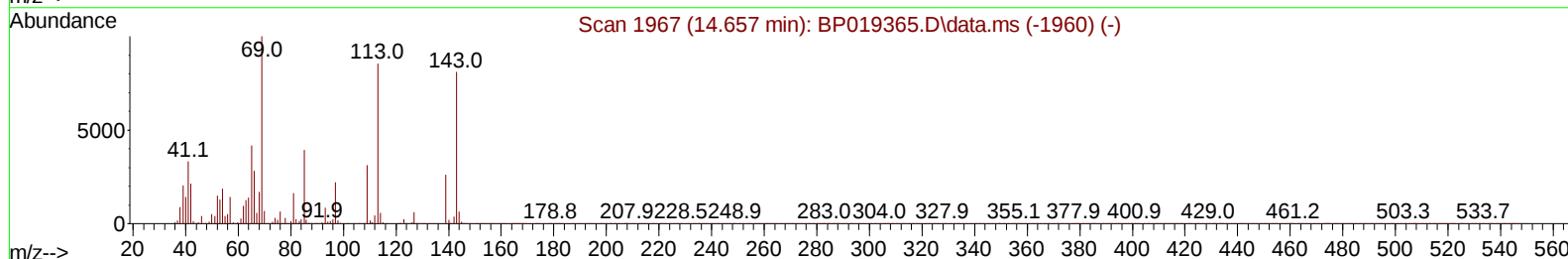
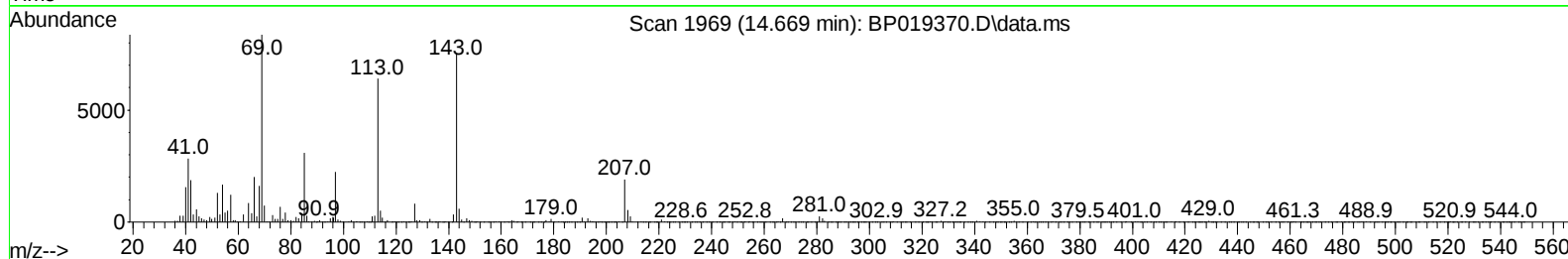
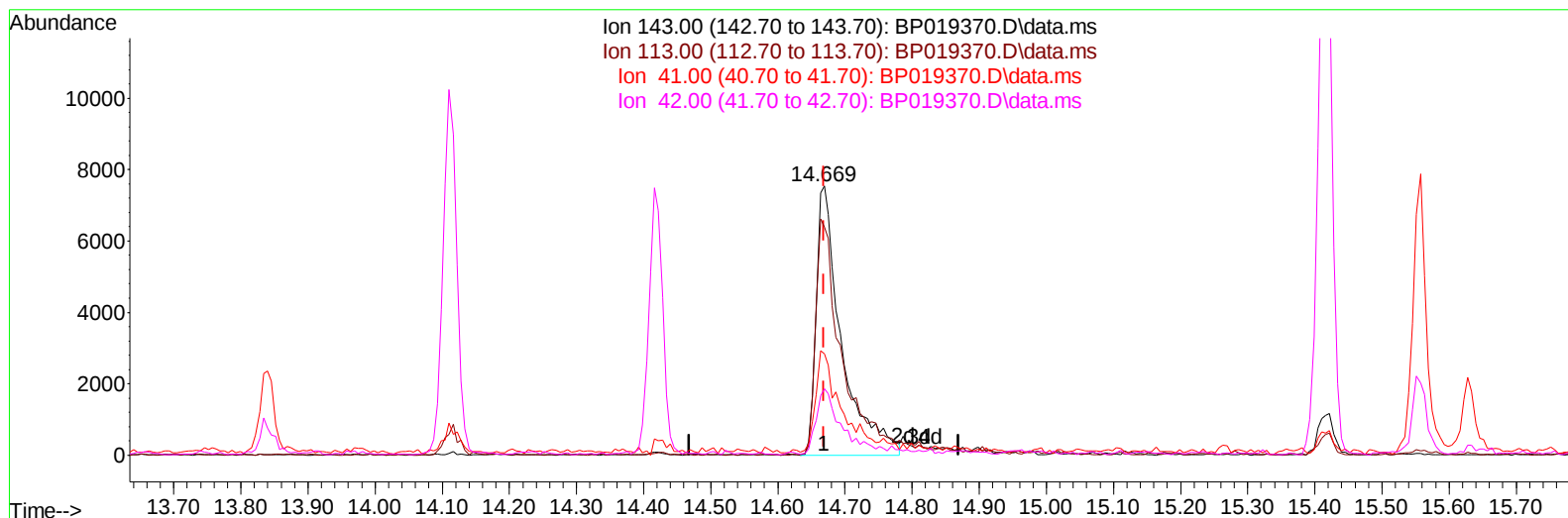
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TIC: BP019370.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 4.35 ng/ul m

response 19752

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	84.85
41.00	39.60	37.71
42.00	25.70	24.66

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Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

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

Quant Time: Jan 16 00:27:57 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	119379	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.569	136	480341	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.422	164	335077	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.175	188	828090	20.000	ng/ul	-0.01
79) Chrysene-d12	21.280	240	807725	20.000	ng/ul	0.00
88) Perylene-d12	23.633	264	807430	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	13182	4.175	ng/uL	0.00
4) Pyridine-d5	3.640	84	76307	8.359	ng/ul	0.00
7) Phenol-d5	6.964	99	65872	5.773	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.117	67	255345	37.415	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.311	132	205833	26.261	ng/ul	0.00
15) 4-Methylphenol-d8	8.499	113	136455	15.075	ng/ul	-0.02
21) Nitrobenzene-d5	8.940	128	154345	38.194	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.663	143	157744	33.595	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.199	165	293809	30.972	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.722	131	371527	31.006	ng/ul	0.00
46) Dimethylphthalate-d6	13.840	166	1113415	37.611	ng/ul	0.00
49) Acenaphthylene-d8	14.110	160	1181043	37.218	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.669	143	19752m	4.345	ng/ul	0.00
60) Fluorene-d10	15.416	176	979643	40.420	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	157799	25.811	ng/ul	0.00
73) Anthracene-d10	17.275	188	1645504	39.842	ng/ul	-0.01
81) Pyrene-d10	19.522	212	2058000	39.074	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.480	264	1788181	40.729	ng/ul	0.00

Target Compounds Qvalue

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

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