

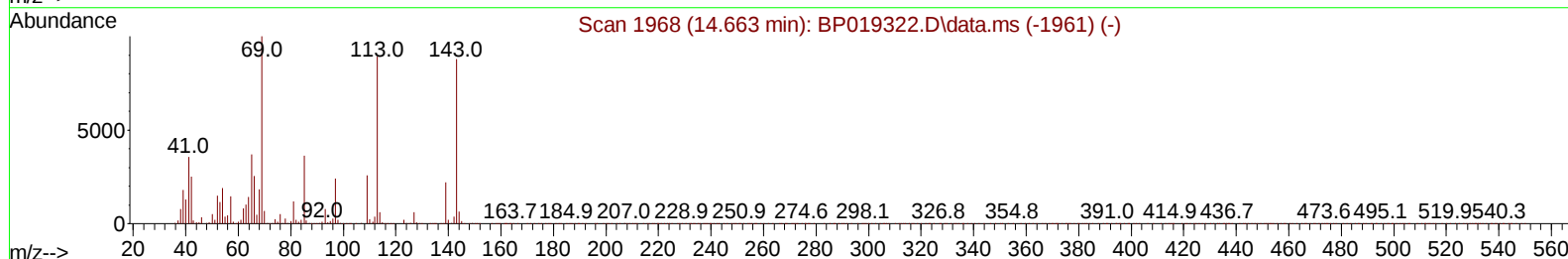
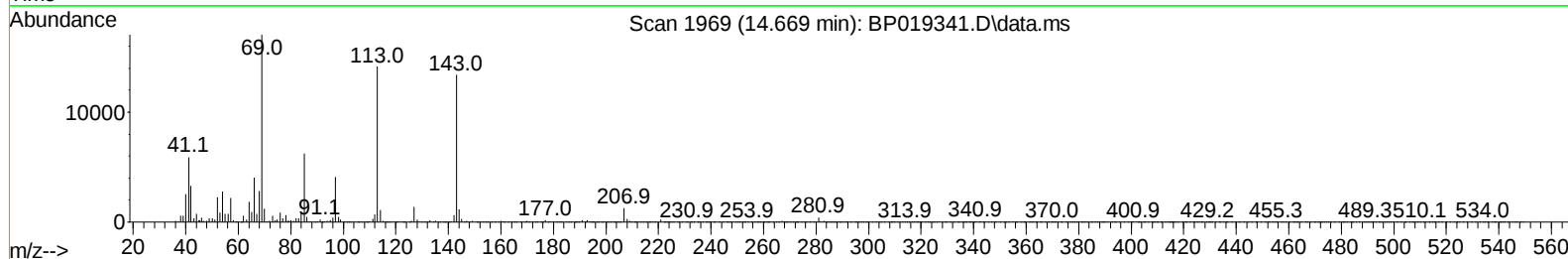
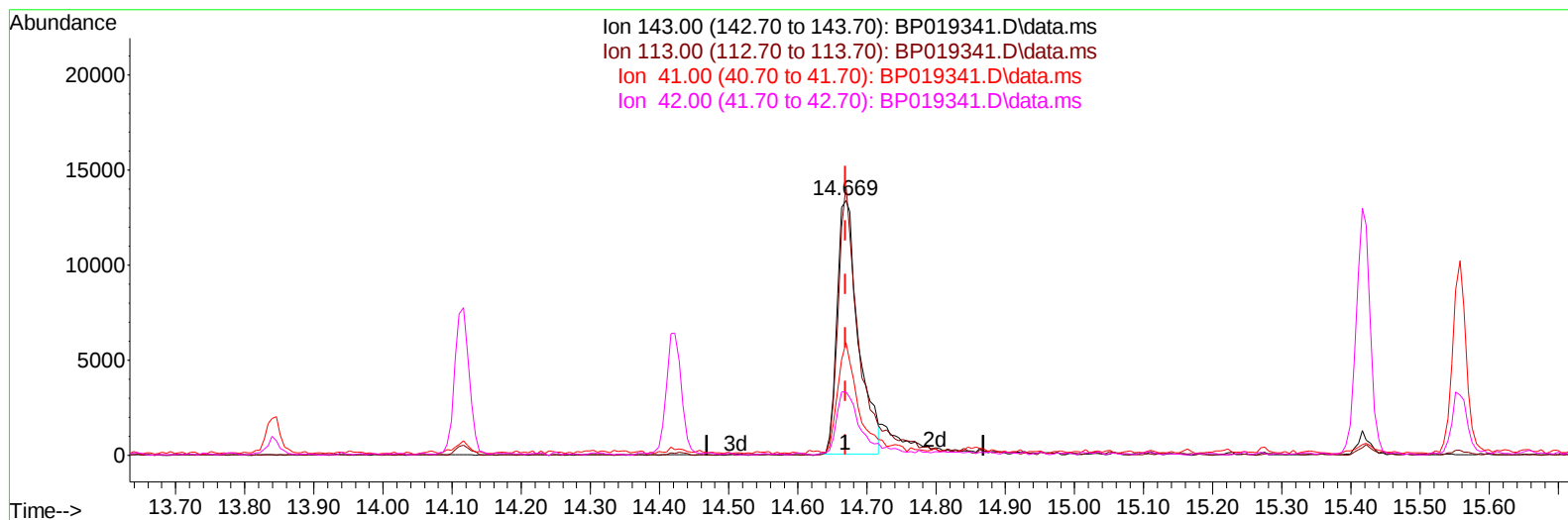
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011224\
Data File : BP019341.D
Acq On : 12 Jan 2024 21:30
Operator : MA/JU
Sample : P1091-14
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0B45

Manual IntegrationsAPPROVED

Quant Time: Jan 12 22:05:22 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 01/15/2024
Supervised By :mohammad ahmed 01/16/2024



TIC: BP019341.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 6.45 ng/u1

response 28520

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	105.61
41.00	39.60	44.15
42.00	25.70	24.80

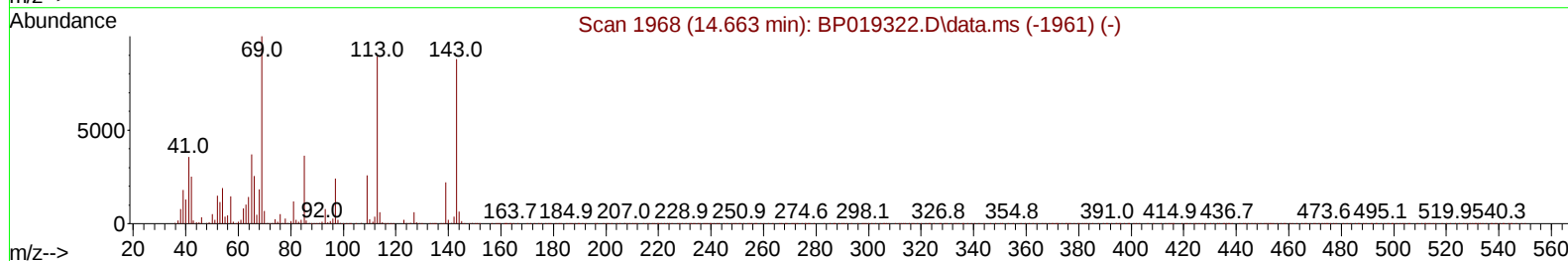
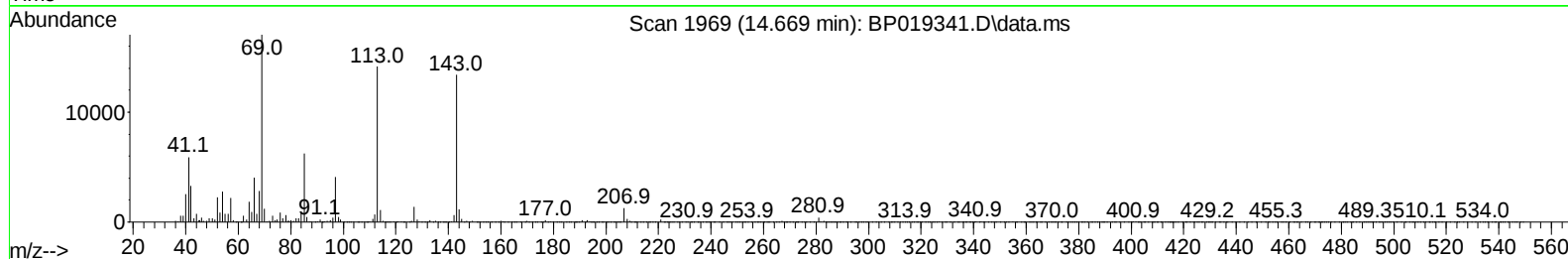
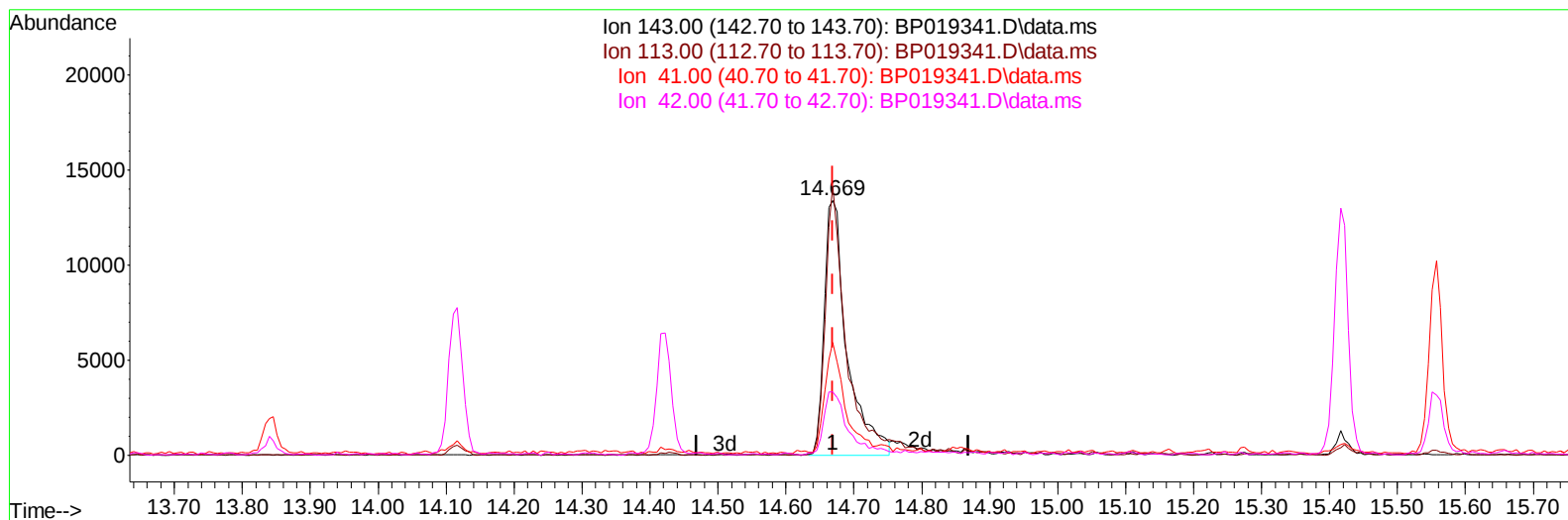
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TIC: BP019341.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.000) 7.04 ng/ul m

response 31159

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	98.20	105.61
41.00	39.60	44.15
42.00	25.70	24.80

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Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 01/15/2024
 Supervised By :mohammad ahmed 01/16/2024

Quant Time: Jan 12 22:06:04 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
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	112928	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.575	136	462560	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.422	164	326164	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.181	188	835061	20.000	ng/ul	0.00
79) Chrysene-d12	21.281	240	895981	20.000	ng/ul	0.00
88) Perylene-d12	23.633	264	964718	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	11644	3.899	ng/uL	0.00
4) Pyridine-d5	3.640	84	78418	9.081	ng/ul	0.00
7) Phenol-d5	6.970	99	64705	5.995	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.122	67	197591	30.606	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	176719	23.835	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	124312	14.518	ng/ul	-0.01
21) Nitrobenzene-d5	8.946	128	120278	30.908	ng/ul	0.00
24) 2-Nitrophenol-d4	9.663	143	136367	30.159	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.205	165	240751	26.355	ng/ul	0.00
31) 4-Chloroaniline-d4	10.728	131	284265	24.635	ng/ul	0.00
46) Dimethylphthalate-d6	13.846	166	923709	32.056	ng/ul	0.00
49) Acenaphthylene-d8	14.116	160	931670	30.162	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	31159m	7.042	ng/ul	0.00
60) Fluorene-d10	15.416	176	818518	34.695	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.557	200	222821	36.142	ng/ul	0.00
73) Anthracene-d10	17.281	188	1624550	39.007	ng/ul	0.00
81) Pyrene-d10	19.528	212	2210134	37.829	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.486	264	2213838	42.203	ng/ul	0.00

Target Compounds Qvalue

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

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