

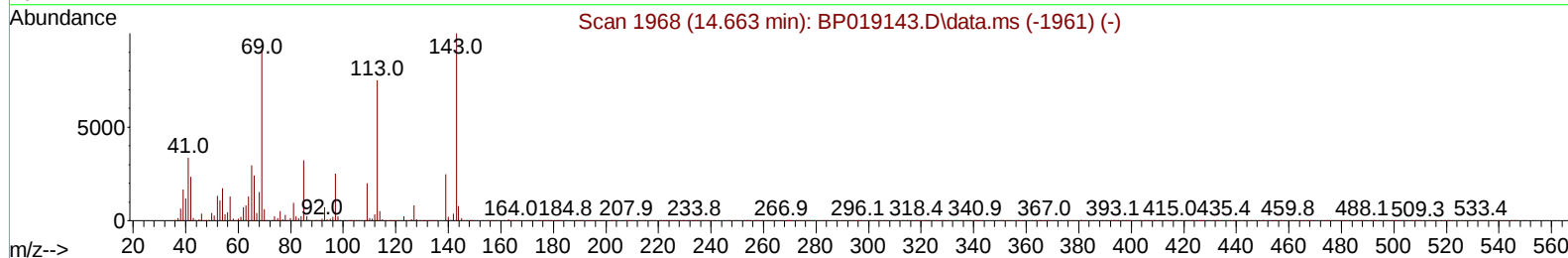
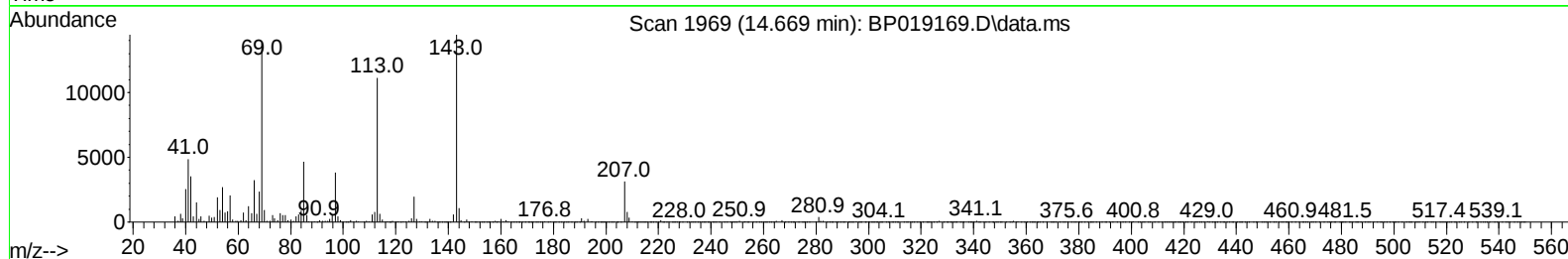
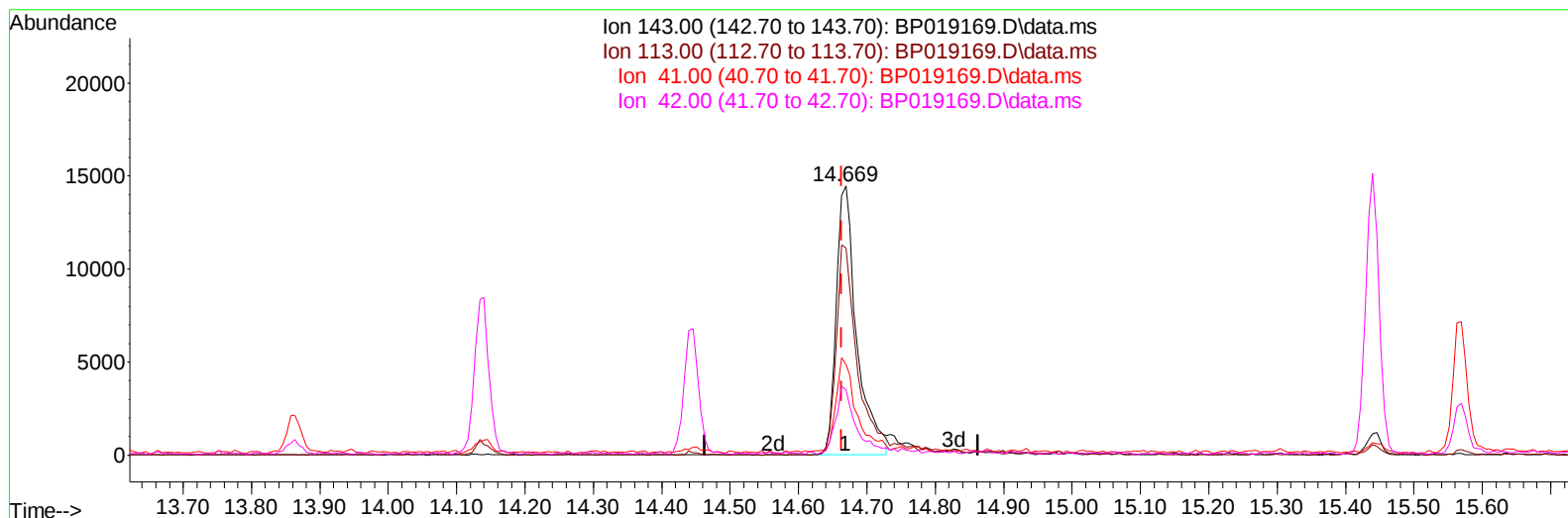
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019169.D
 Acq On : 29 Dec 2023 16:57
 Operator : MA/JU
 Sample : 06020-07
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7W8

Manual IntegrationsAPPROVED

Quant Time: Dec 29 18:20:51 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 01/01/2024
 Supervised By :mohammad ahmed 01/01/2024



TIC: BP019169.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.006) 4.91 ng/ul

response 30154

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	77.10
41.00	33.40	33.58
42.00	23.00	24.28

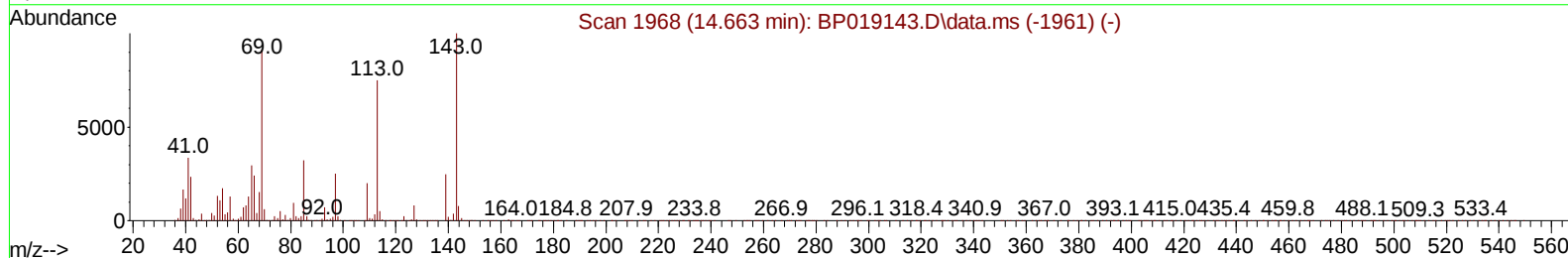
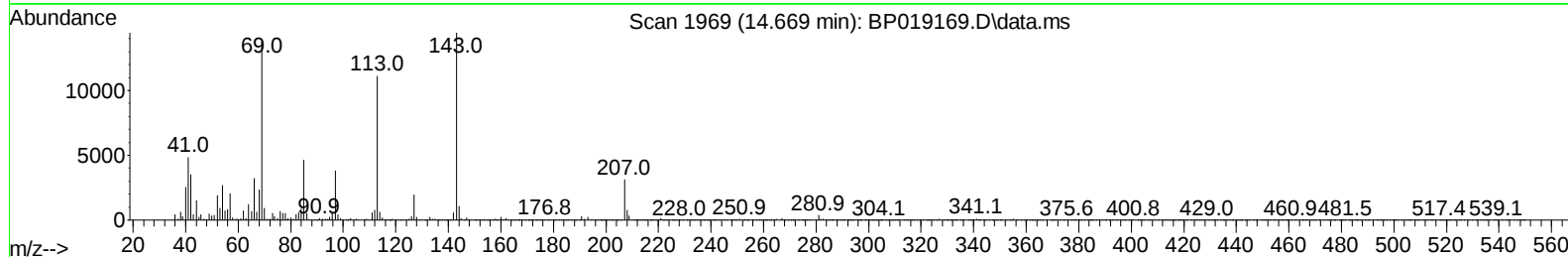
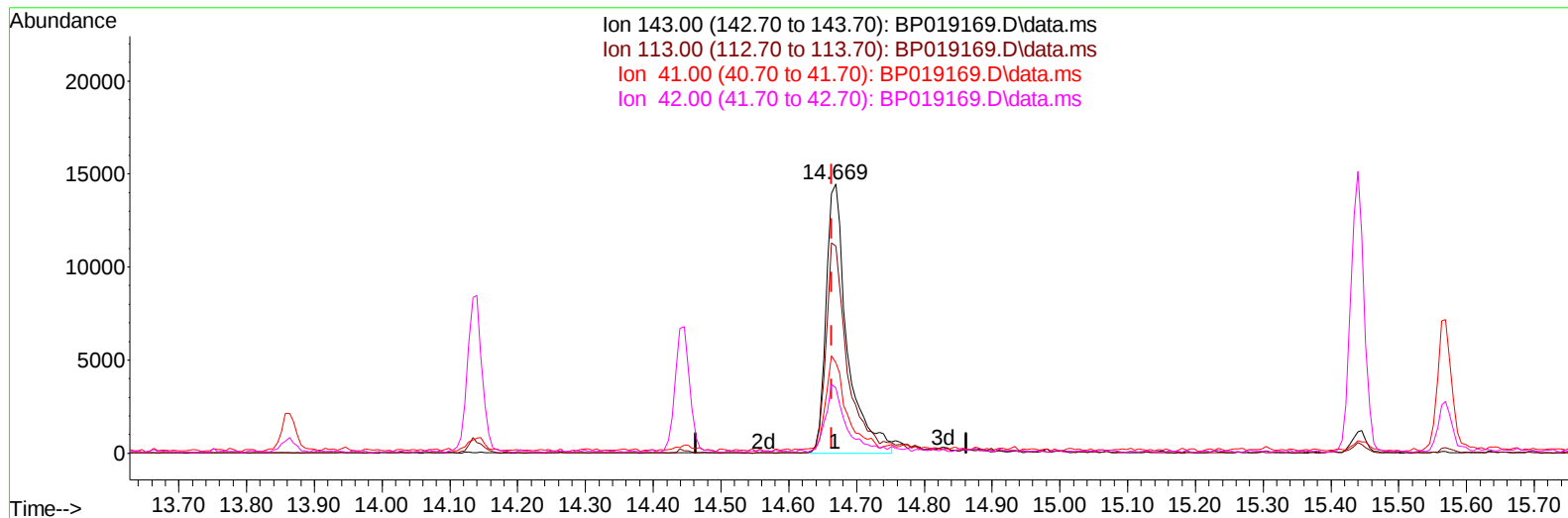
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TIC: BP019169.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.006) 5.12 ng/ul m

response 31451

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	77.10
41.00	33.40	33.58
42.00	23.00	24.28

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 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 01/01/2024
 Supervised By :mohammad ahmed 01/01/2024

Quant Time: Dec 30 01:30:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 29 03:48:11 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	193057	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136	700445	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	411521	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	935328	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1012120	20.000	ng/ul	0.00
88) Perylene-d12	23.669	264	1215308	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	23599	4.175	ng/uL	0.00
4) Pyridine-d5	3.658	84	132920	9.368	ng/ul	0.00
7) Phenol-d5	6.975	99	114177	6.595	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	303078	30.260	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	319337	22.880	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	200045	14.841	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	179477	29.641	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	179241	27.391	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	346096	26.192	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	441928	27.190	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1200118	33.017	ng/ul	0.00
49) Acenaphthylene-d8	14.140	160	1260422	31.958	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	31451m	5.124	ng/ul	0.00
60) Fluorene-d10	15.440	176	1076412	34.278	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	163284	26.629	ng/ul	0.00
73) Anthracene-d10	17.298	188	1836416	37.596	ng/ul	0.00
81) Pyrene-d10	19.545	212	2297966	37.255	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.516	264	2644496	37.856	ng/ul	0.00

Target Compounds Qvalue

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

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