

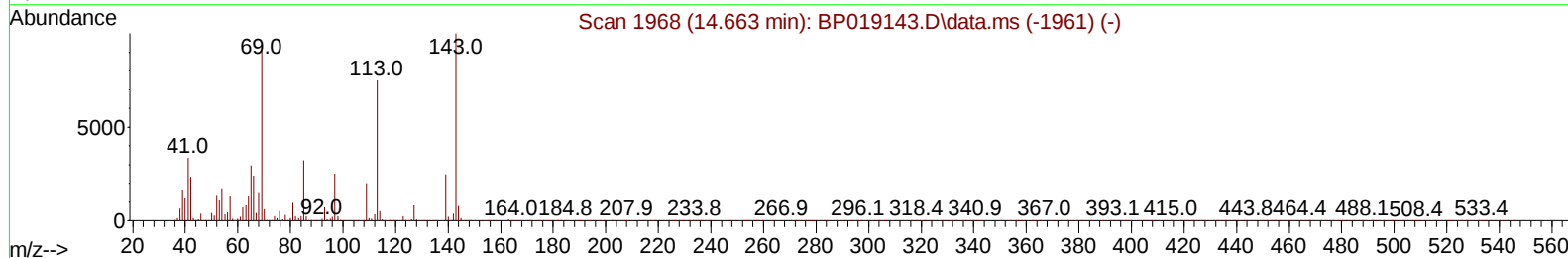
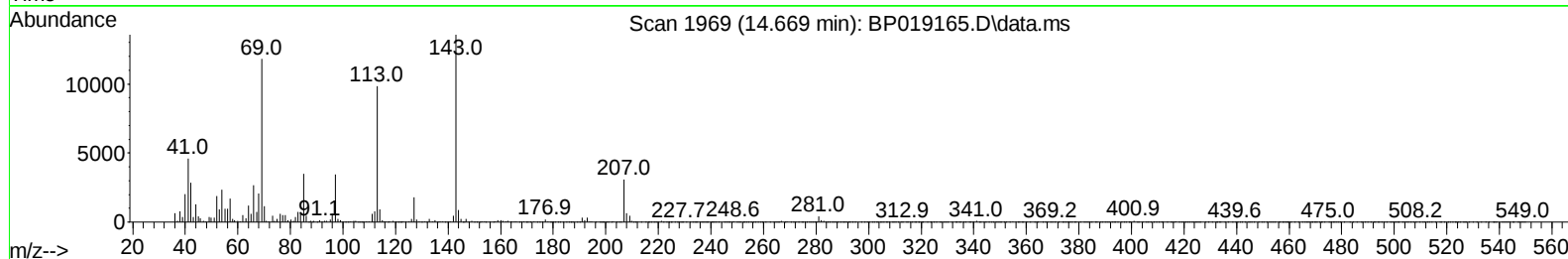
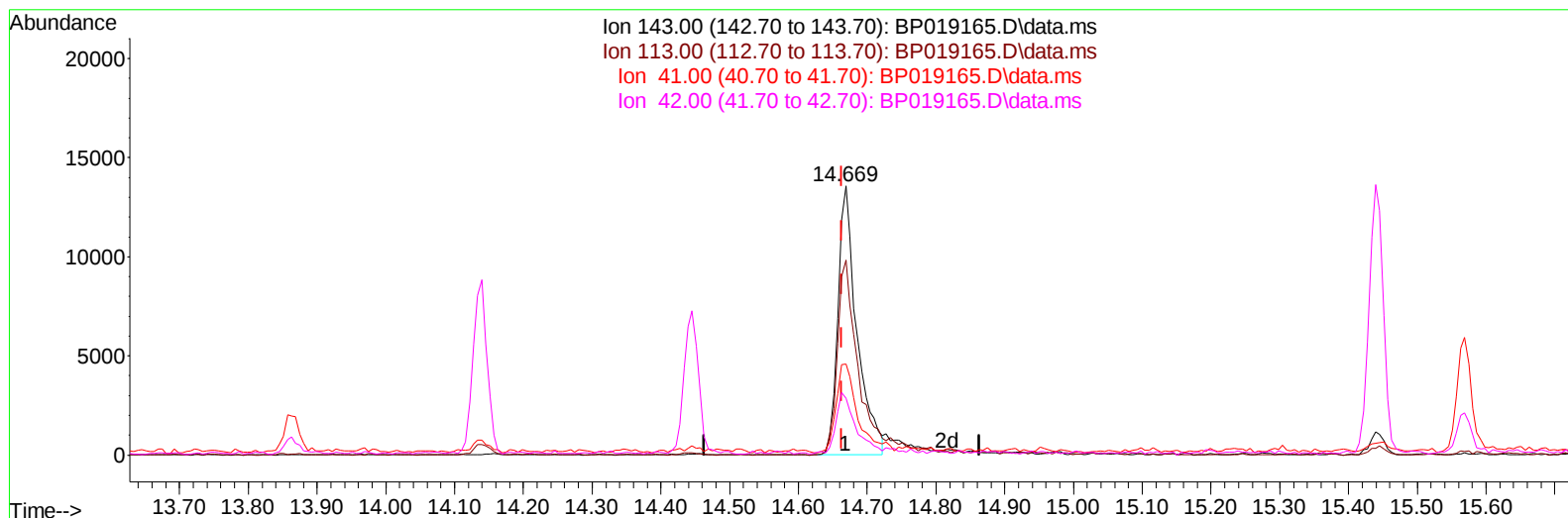
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019165.D
 Acq On : 29 Dec 2023 14:41
 Operator : MA/JU
 Sample : 06021-08
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH7Z0

Manual IntegrationsAPPROVED

Quant Time: Dec 29 16:21:41 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: BP019165.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.006) 4.25 ng/u1

response 26013

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	72.42
41.00	33.40	33.90
42.00	23.00	21.16

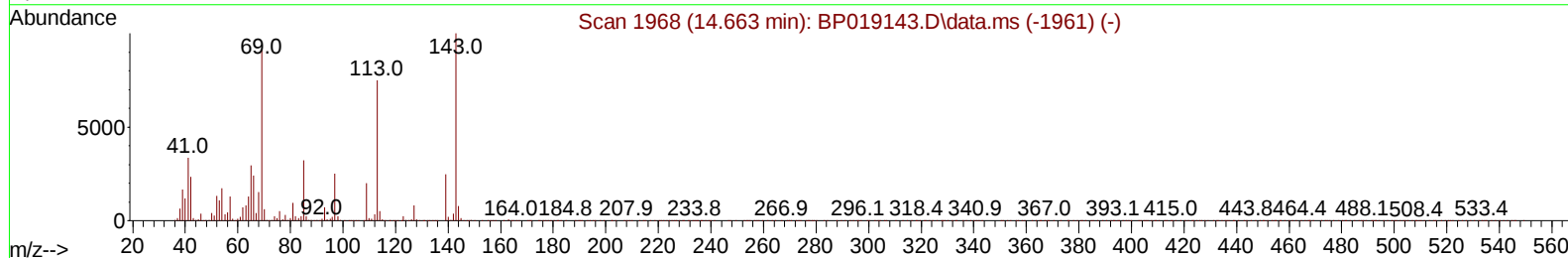
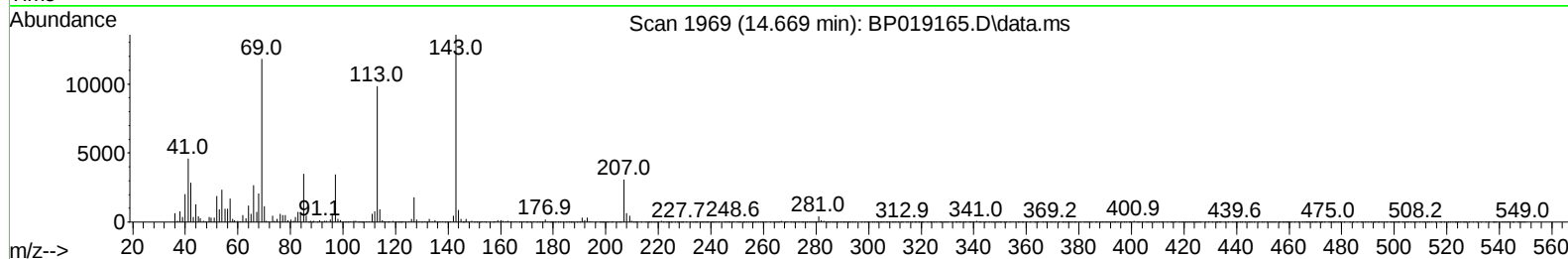
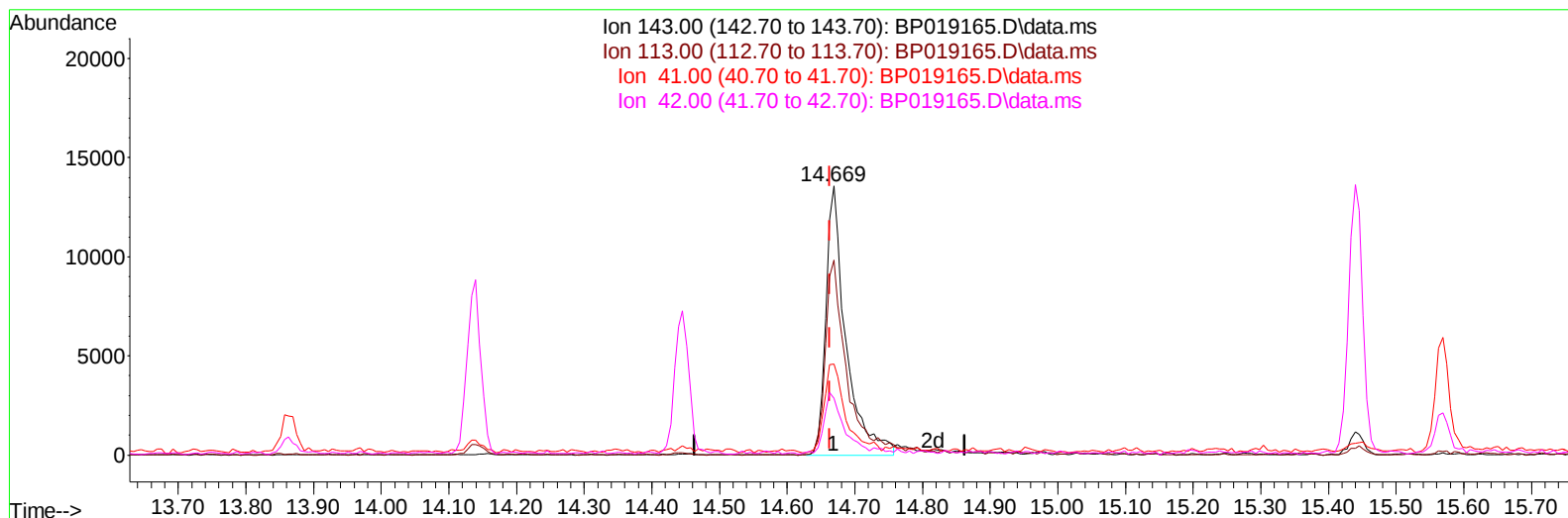
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TIC: BP019165.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.669min (+ 0.006) 4.51 ng/ul m

response 27646

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	72.42
41.00	33.40	33.90
42.00	23.00	21.16

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 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
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Manual IntegrationsAPPROVED

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

Quant Time: Dec 30 01:24:07 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	7.805	152	183773	20.000	ng/ul	0.00
20) Naphthalene-d8	10.598	136	681707	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	410655	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	949256	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1039970	20.000	ng/ul	0.00
88) Perylene-d12	23.668	264	1256920	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	22782	4.234	ng/uL	0.00
4) Pyridine-d5	3.658	84	80541	5.963	ng/ul	0.00
7) Phenol-d5	6.981	99	111568	6.770	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	294660	30.906	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	302989	22.805	ng/ul	0.00
15) 4-Methylphenol-d8	8.522	113	195511	15.238	ng/ul	0.00
21) Nitrobenzene-d5	8.969	128	165375	28.063	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	160953	25.272	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	340603	26.484	ng/ul	0.00
31) 4-Chloroaniline-d4	10.746	131	358217	22.645	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1127723	31.090	ng/ul	0.00
49) Acenaphthylene-d8	14.139	160	1245745	31.652	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	27646m	4.514	ng/ul	0.00
60) Fluorene-d10	15.439	176	996155	31.789	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	126140	20.270	ng/ul	0.00
73) Anthracene-d10	17.298	188	1663095	33.548	ng/ul	0.00
81) Pyrene-d10	19.545	212	2112162	33.325	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.521	264	2442961	33.814	ng/ul	0.00

Target Compounds Qvalue

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

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