

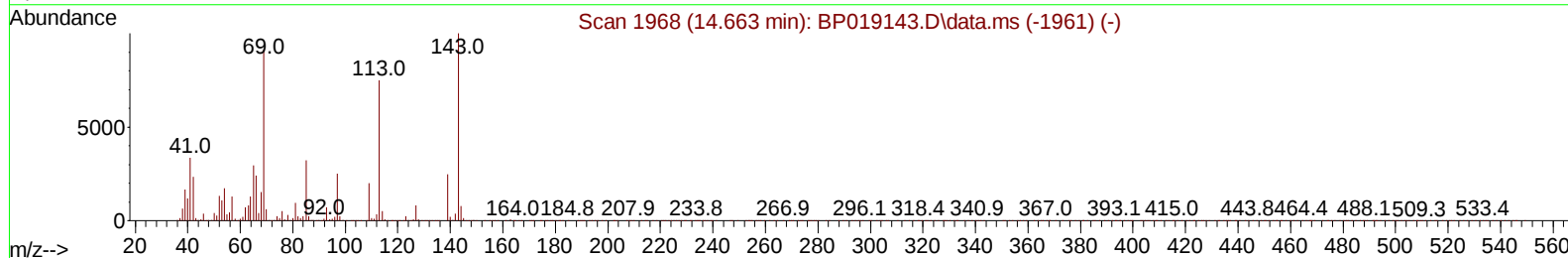
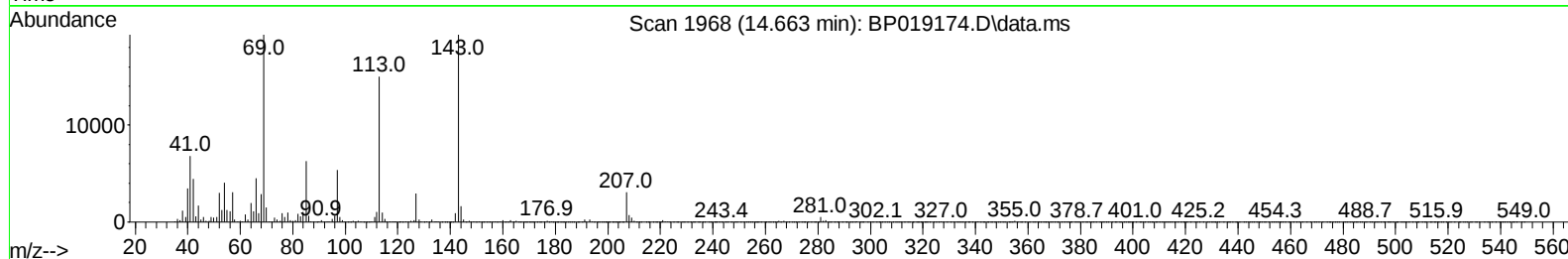
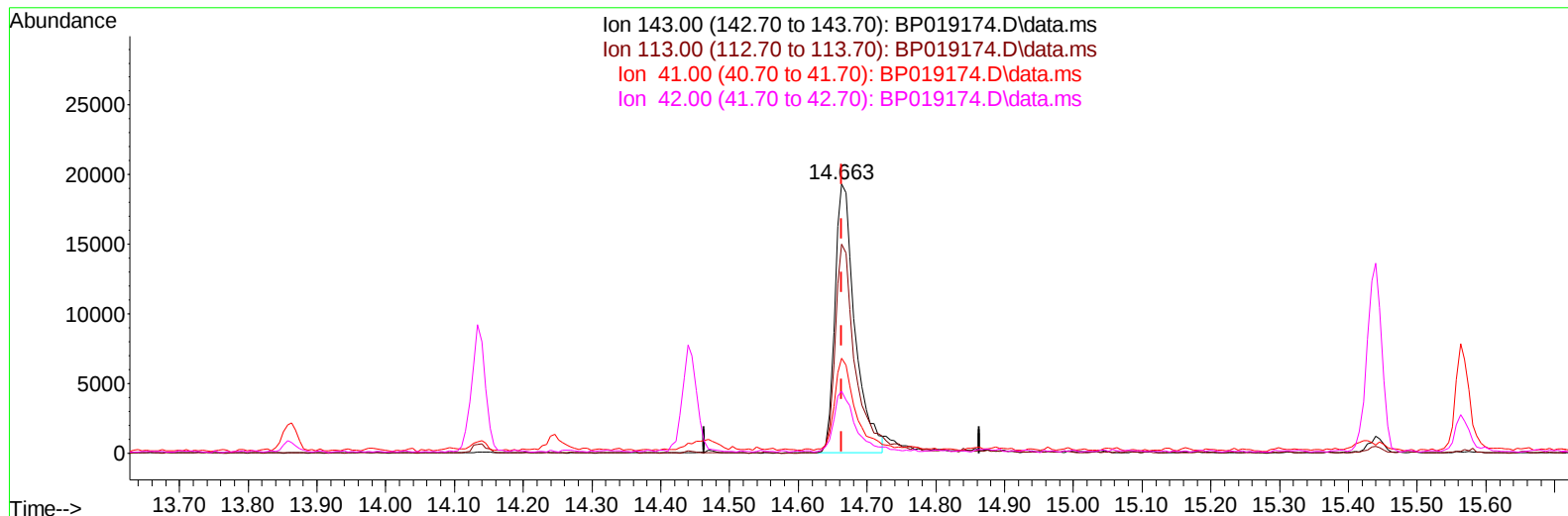
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
Data File : BP019174.D
Acq On : 29 Dec 2023 19:48
Operator : MA/JU
Sample : 06021-02
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH7Y4

Manual IntegrationsAPPROVED

Quant Time: Dec 29 22:01:25 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: BP019174.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.663min (-0.000) 6.13 ng/u1

response 39270

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	77.63
41.00	33.40	35.34
42.00	23.00	23.12

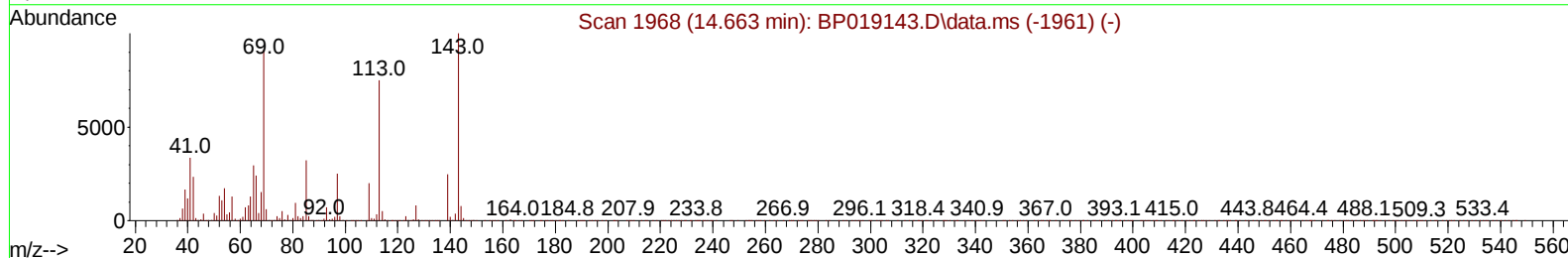
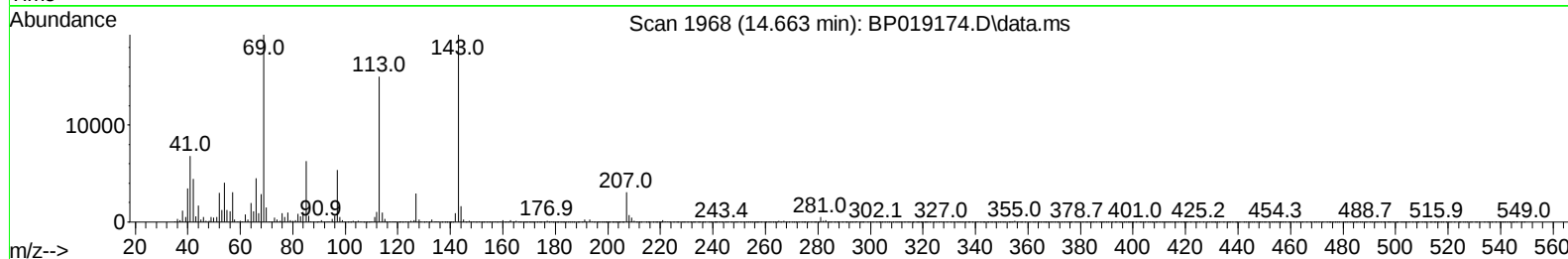
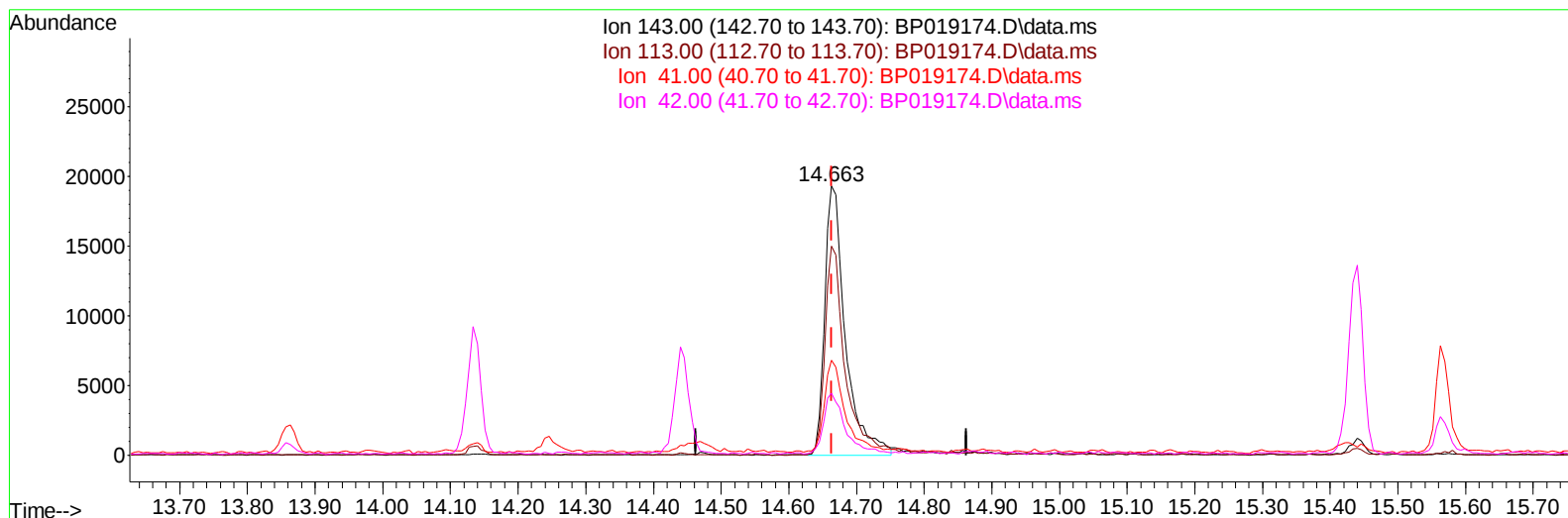
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TIC: BP019174.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.663min (-0.000) 6.37 ng/ul m

response 40798

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	77.63
41.00	33.40	35.34
42.00	23.00	23.12

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Reviewed By :Yogesh Patel 01/01/2024
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Quant Time: Dec 29 22:02:16 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	182309	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	707338	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.445	164	429447	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	934706	20.000	ng/ul	0.00
79) Chrysene-d12	21.304	240	1045642	20.000	ng/ul	0.00
88) Perylene-d12	23.663	264	1245920	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	23836	4.465	ng/uL	0.00
4) Pyridine-d5	3.652	84	129677	9.679	ng/ul	0.00
7) Phenol-d5	6.975	99	133148	8.144	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	299717	31.689	ng/ul	0.00
11) 2-Chlorophenol-d4	7.334	132	322868	24.496	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	221946	17.437	ng/ul	0.00
21) Nitrobenzene-d5	8.963	128	180405	29.504	ng/ul	0.00
24) 2-Nitrophenol-d4	9.687	143	190183	28.780	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	367028	27.505	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	386652	23.557	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1200062	31.637	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	1247816	30.317	ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	40798m	6.369	ng/ul	0.00
60) Fluorene-d10	15.440	176	1015572	30.991	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.569	200	159548	26.037	ng/ul	0.00
73) Anthracene-d10	17.298	188	1673048	34.274	ng/ul	0.00
81) Pyrene-d10	19.545	212	2188897	34.349	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.510	264	2464960	34.419	ng/ul	0.00

Target Compounds Qvalue

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

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