

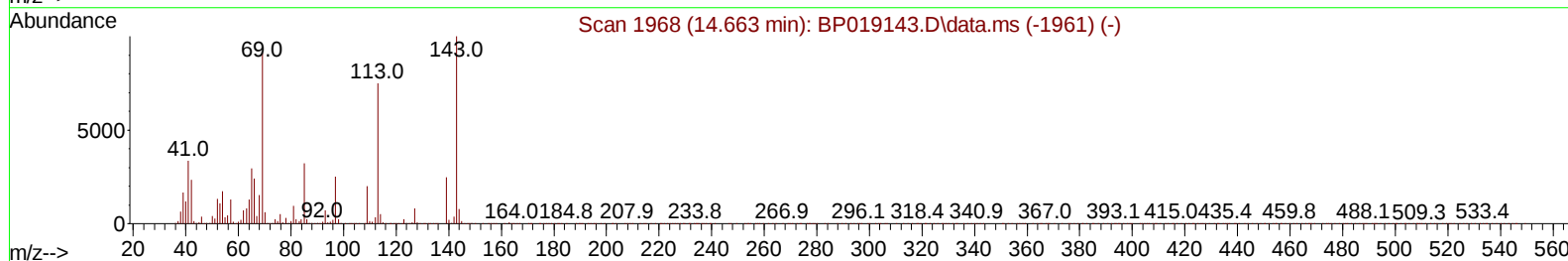
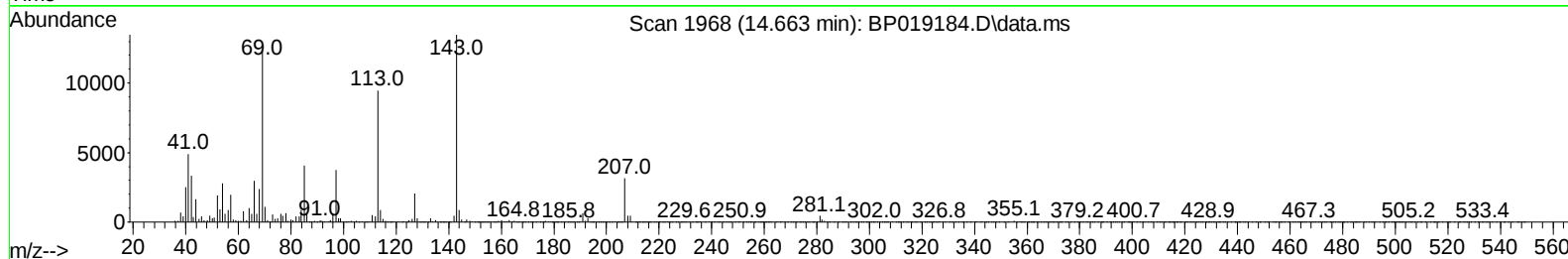
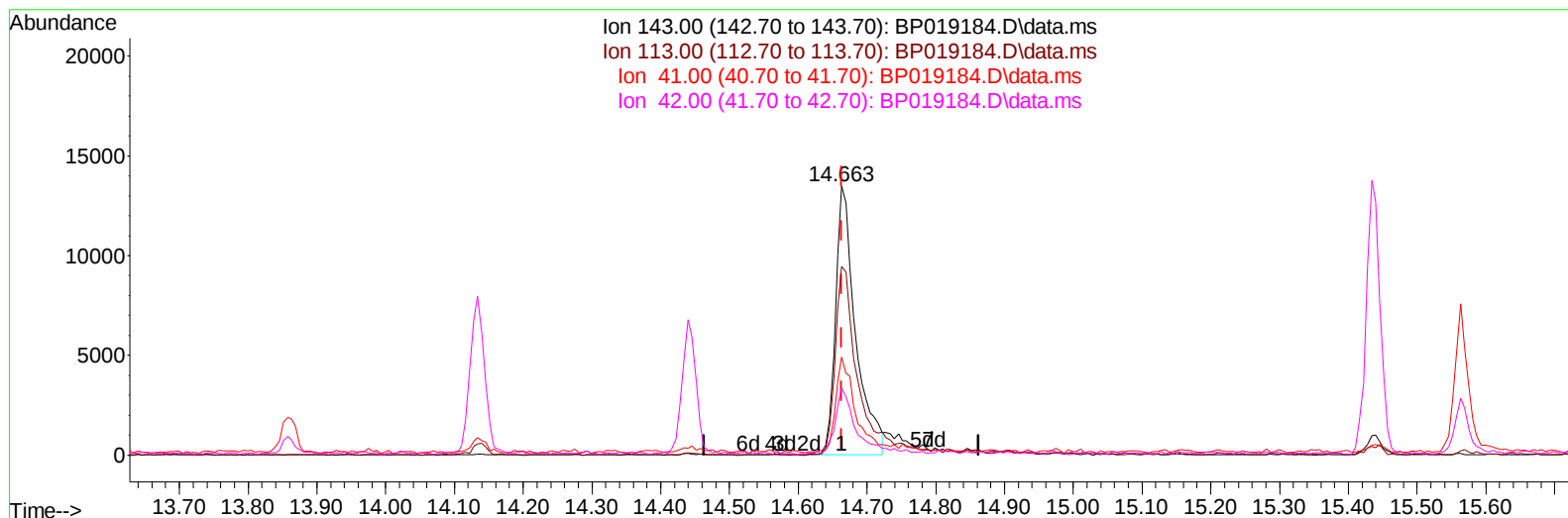
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122823\
 Data File : BP019184.D
 Acq On : 30 Dec 2023 02:05
 Operator : MA/JU
 Sample : 06020-01
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH3N9

Manual IntegrationsAPPROVED

Quant Time: Dec 30 03:35:41 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: BP019184.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.663min (+ 0.000) 5.18 ng/u1

response 27426

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	70.17
41.00	33.40	36.45
42.00	23.00	24.84

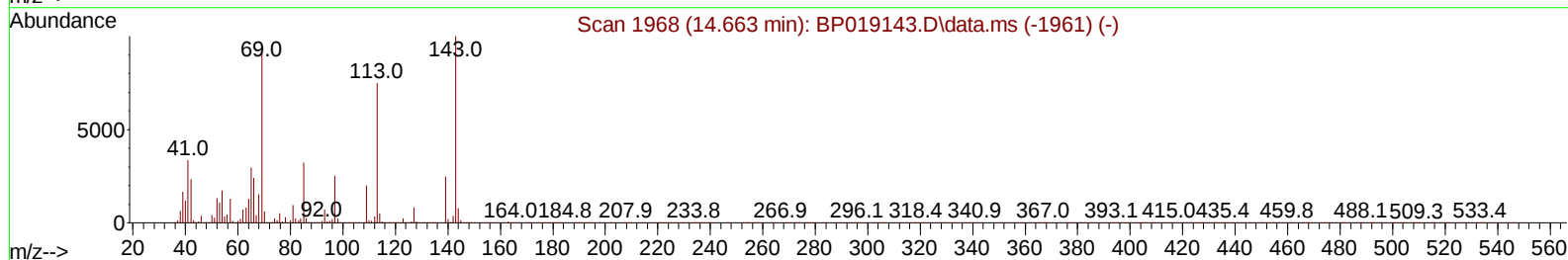
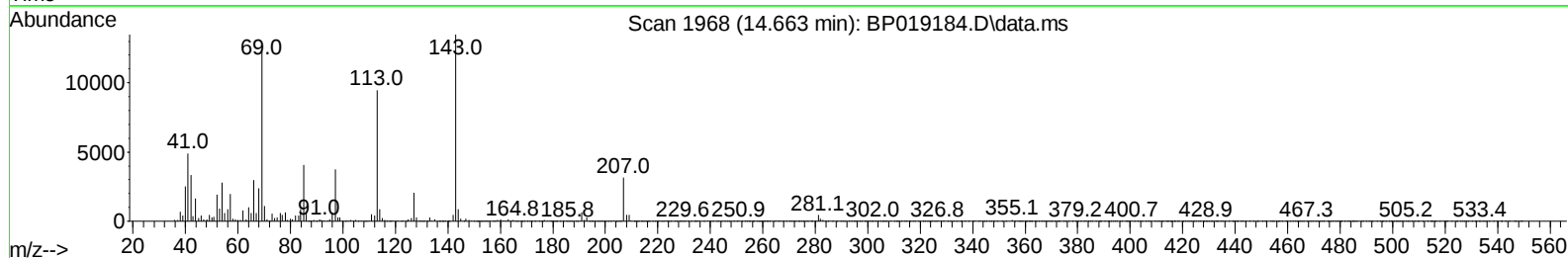
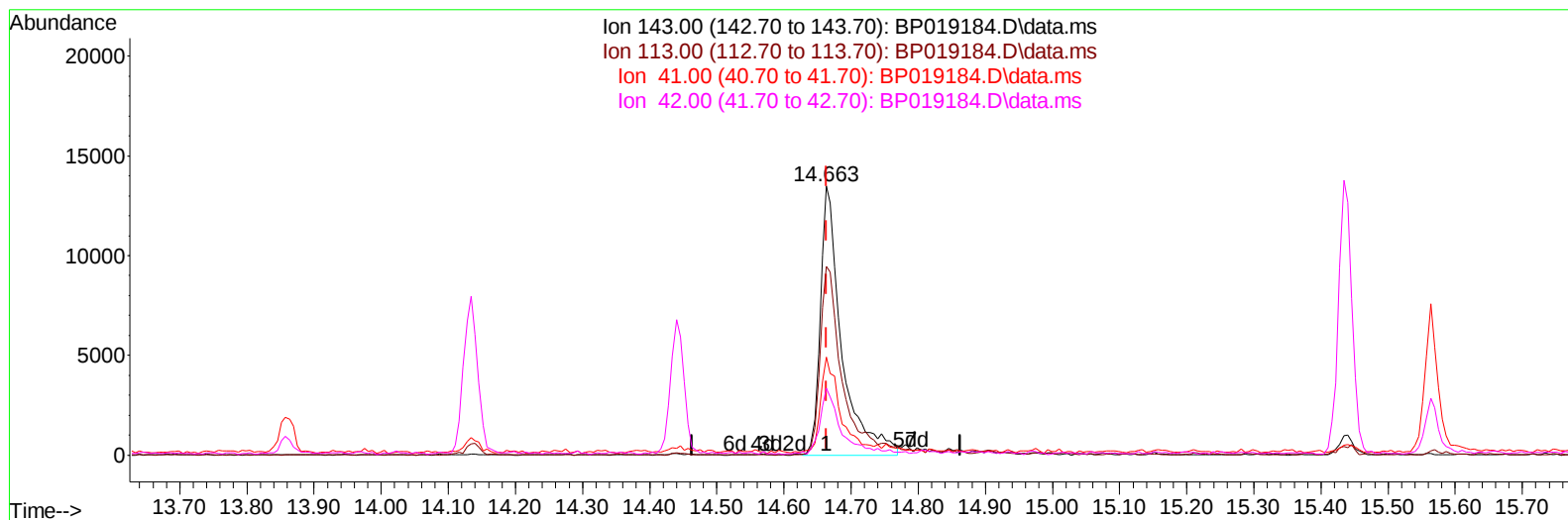
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TIC: BP019184.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.663min (+ 0.000) 5.62 ng/ul m

response 29715

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	73.50	70.17
41.00	33.40	36.45
42.00	23.00	24.84

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 01/01/2024
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Quant Time: Dec 30 03:37:56 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	175393	20.000	ng/ul	0.00
20) Naphthalene-d8	10.593	136	632838	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.440	164	354711	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.198	188	784067	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	868401	20.000	ng/ul	-0.01
88) Perylene-d12	23.663	264	1052163	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.234	96	25022	4.872	ng/uL	0.00
4) Pyridine-d5	3.658	84	124954	9.694	ng/ul	0.00
7) Phenol-d5	6.975	99	114890	7.304	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	304900	33.508	ng/ul	0.00
11) 2-Chlorophenol-d4	7.328	132	316389	24.951	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	184810	15.092	ng/ul	0.00
21) Nitrobenzene-d5	8.964	128	178026	32.543	ng/ul	0.00
24) 2-Nitrophenol-d4	9.681	143	178463	30.185	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.222	165	331157	27.738	ng/ul	0.00
31) 4-Chloroaniline-d4	10.740	131	403648	27.488	ng/ul	0.00
46) Dimethylphthalate-d6	13.863	166	1050011	33.514	ng/ul	0.00
49) Acenaphthylene-d8	14.134	160	1151775	33.880	ng/ul	0.00
54) 4-Nitrophenol-d4	14.663	143	29715m	5.617	ng/ul	0.00
60) Fluorene-d10	15.440	176	931215	34.404	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.563	200	147208	28.639	ng/ul	-0.01
73) Anthracene-d10	17.298	188	1587392	38.768	ng/ul	0.00
81) Pyrene-d10	19.545	212	2007499	37.932	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.510	264	2388334	39.491	ng/ul	0.00

Target Compounds Qvalue

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

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